

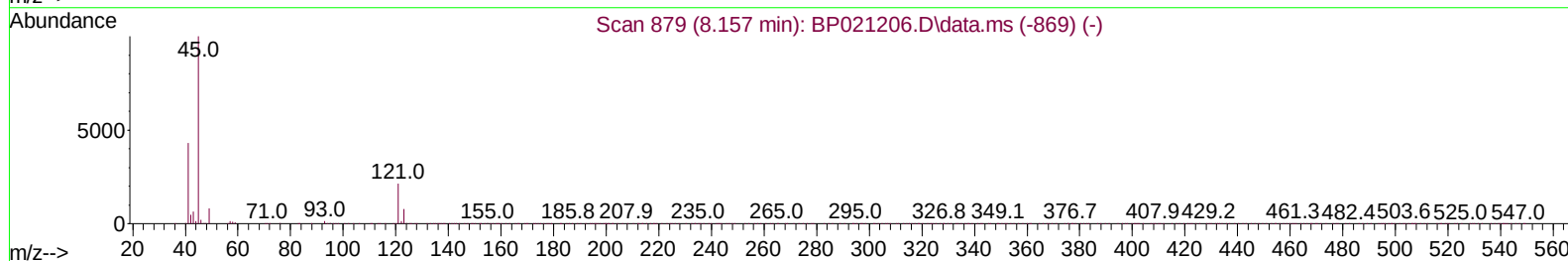
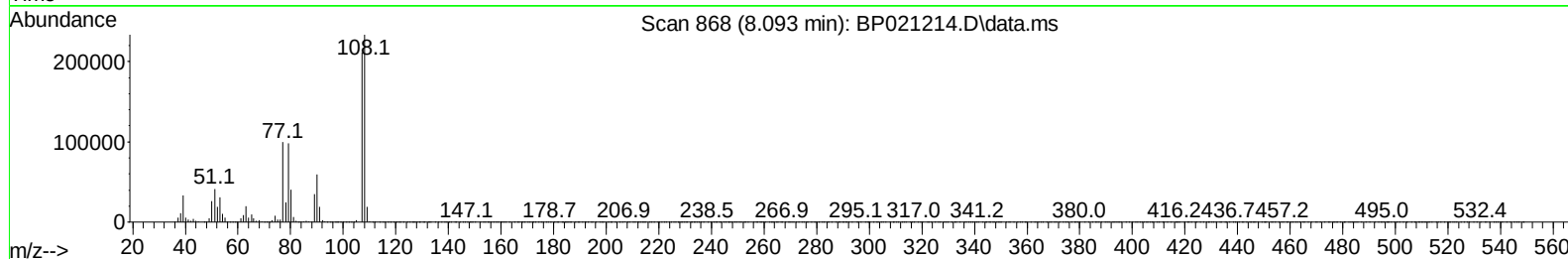
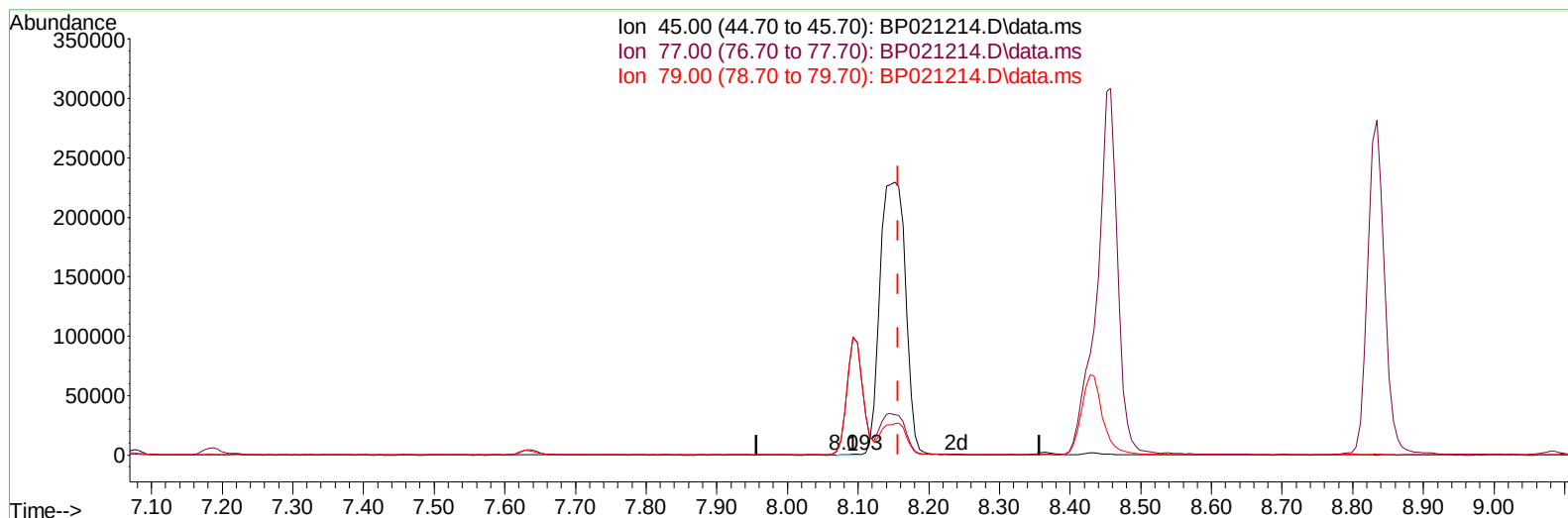
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021214.D
Acq On : 29 Jul 2024 17:50
Operator : MA/JU
Sample : PB162189BS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SLCS189

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/30/2024
Supervised By :mohammad ahmed 09/04/2024

Quant Time: Jul 29 19:08:47 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration



TIC: BP021214.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.093min (-0.064) 0.03 ng/u1

response 732

Ion	Exp%	Act%
45.00	100.00	100.00
77.00	13.60	15462.52#
79.00	10.10	15272.16#
0.00	0.00	0.00

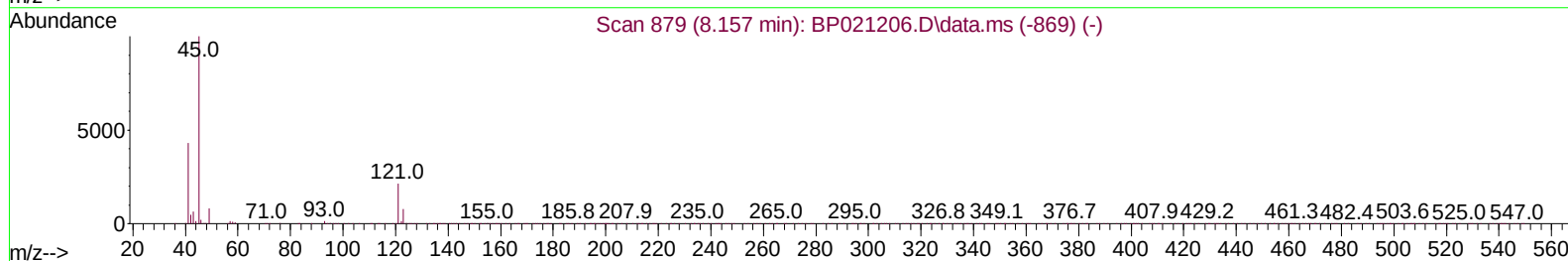
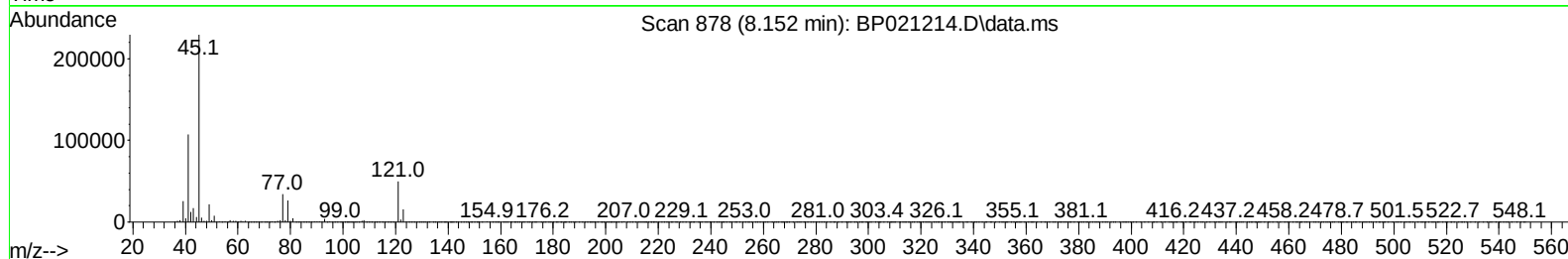
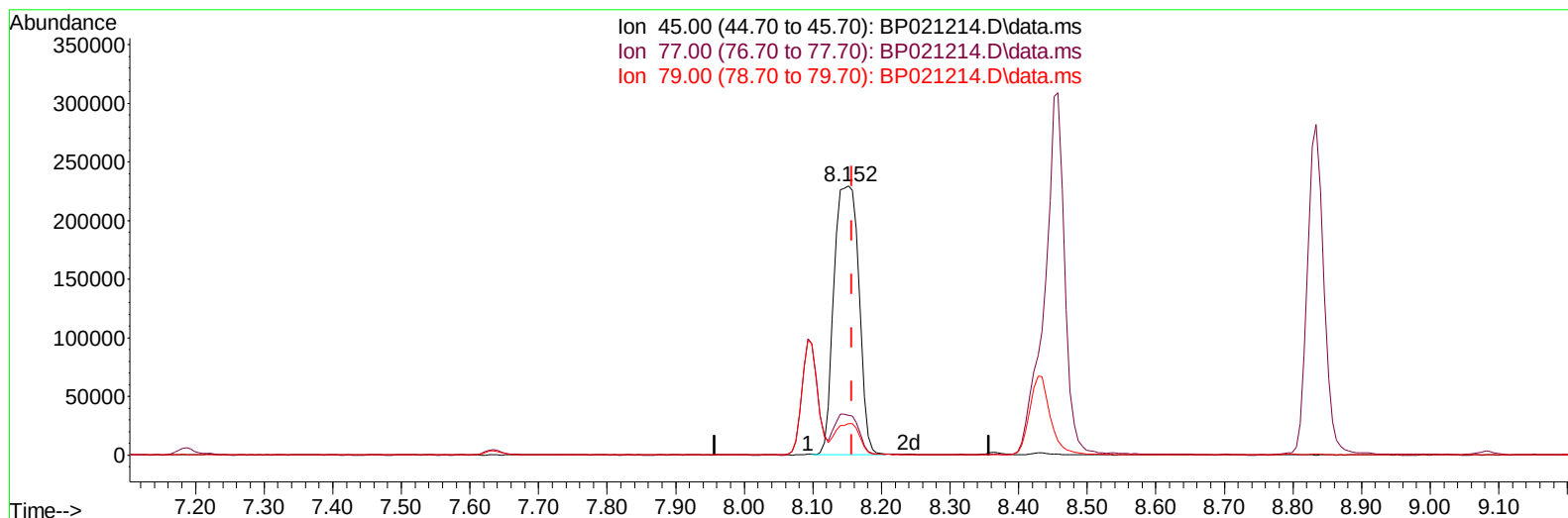
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TIC: BP021214.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.152min (-0.005) 26.73 ng/ul m

response 580226

Ion	Exp%	Act%
45.00	100.00	100.00
77.00	13.60	14.78
79.00	10.10	11.62
0.00	0.00	0.00

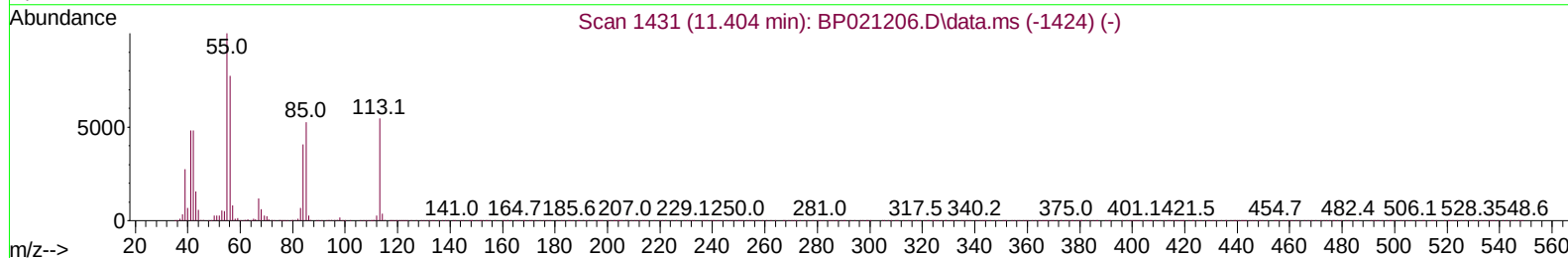
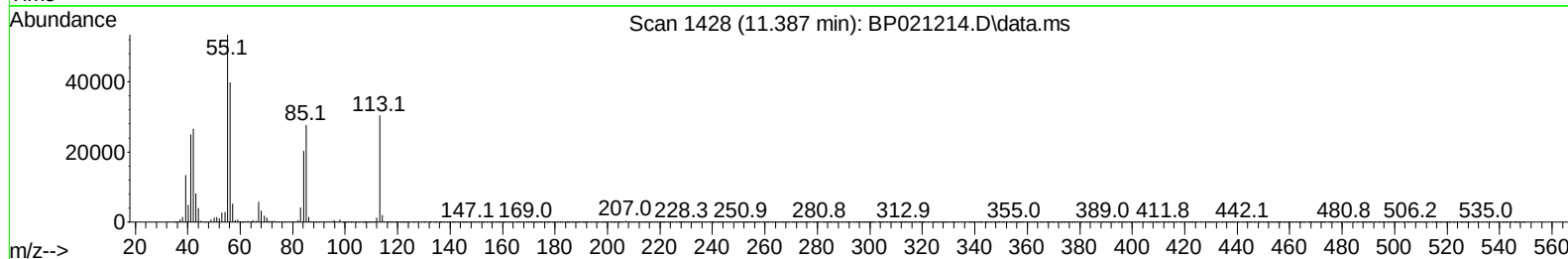
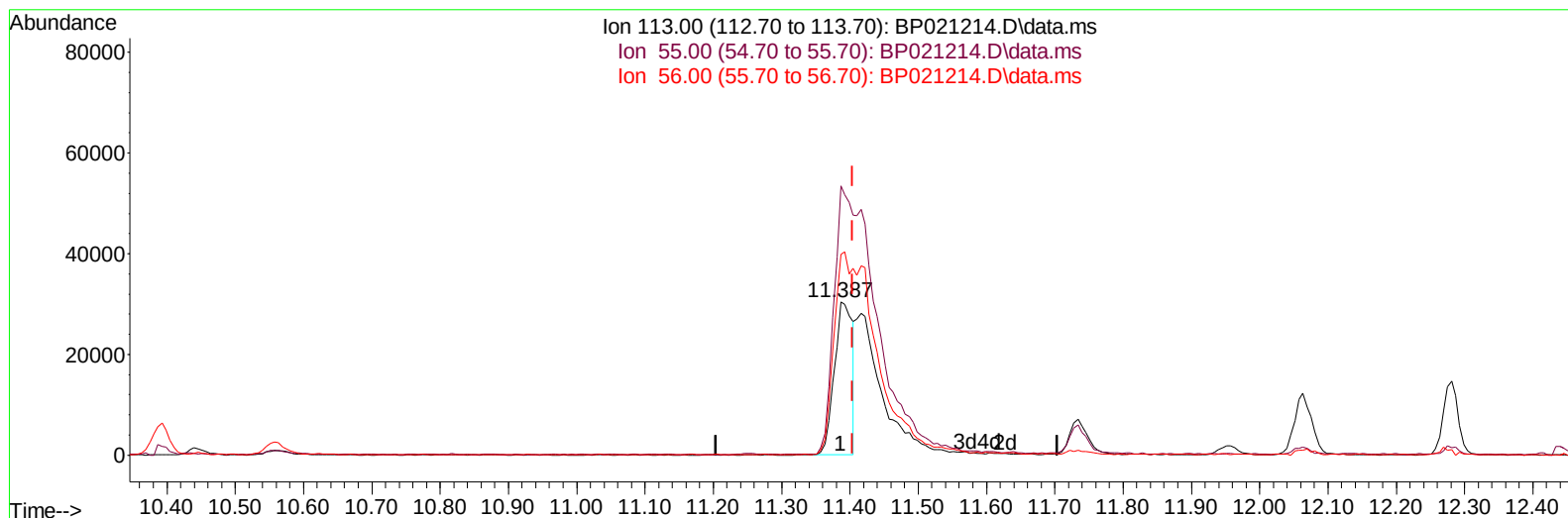
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
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 ALS Vial : 11 Sample Multiplier: 1

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TIC: BP021214.D\data.ms

(34) Caprolactam

11.387min (-0.017) 13.37 ng/ul

response 56373

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	175.30	175.90
56.00	131.90	131.39
0.00	0.00	0.00

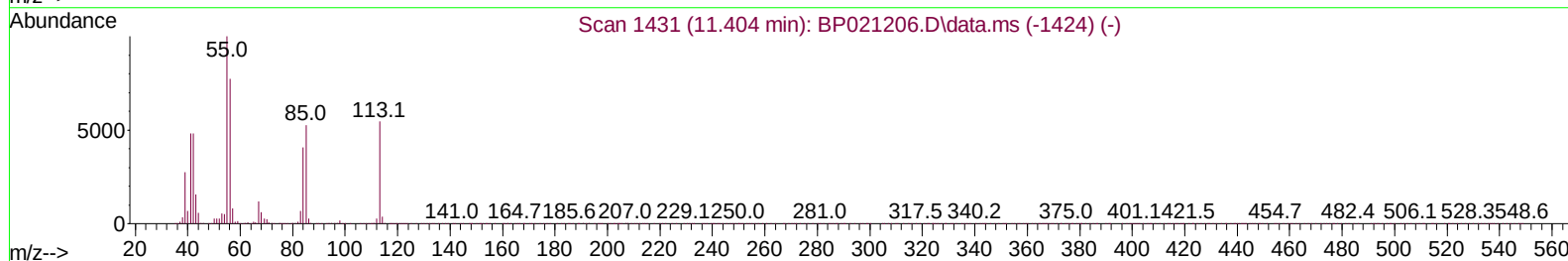
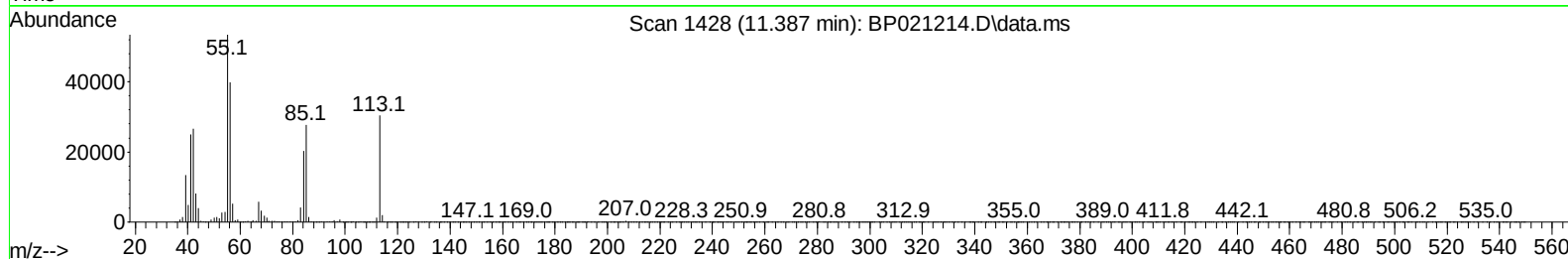
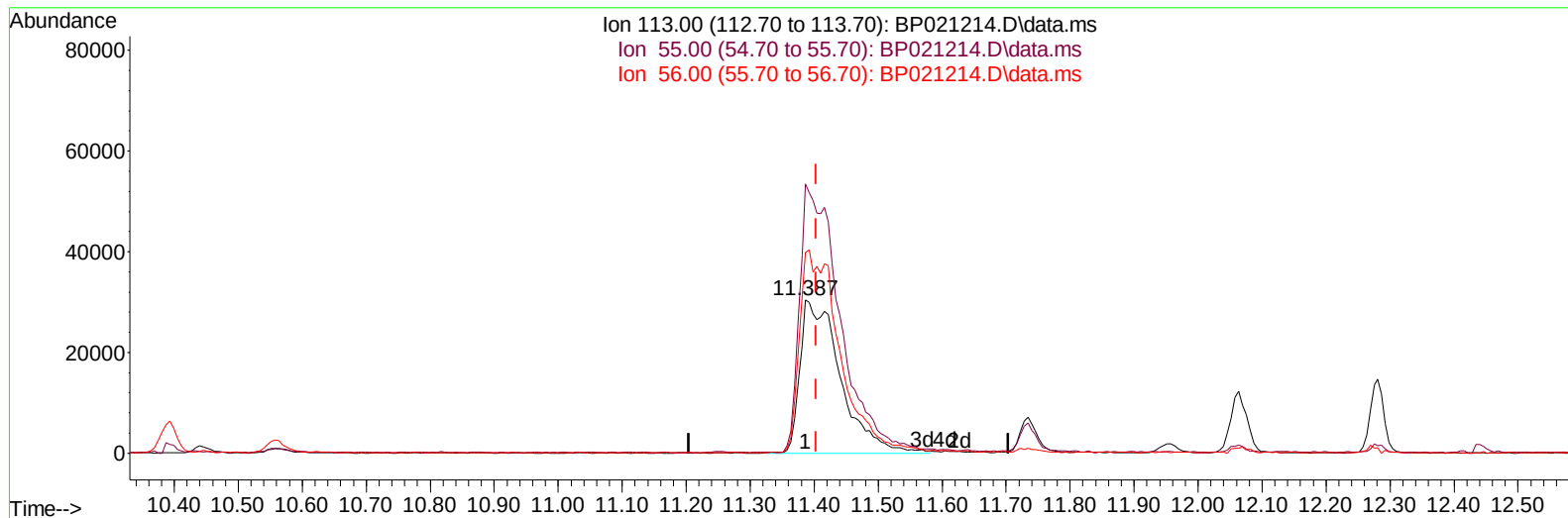
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
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TIC: BP021214.D\data.ms

(34) Caprolactam

11.387min (-0.017) 31.59 ng/ul m

response 133213

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	175.30	175.90
56.00	131.90	131.39
0.00	0.00	0.00

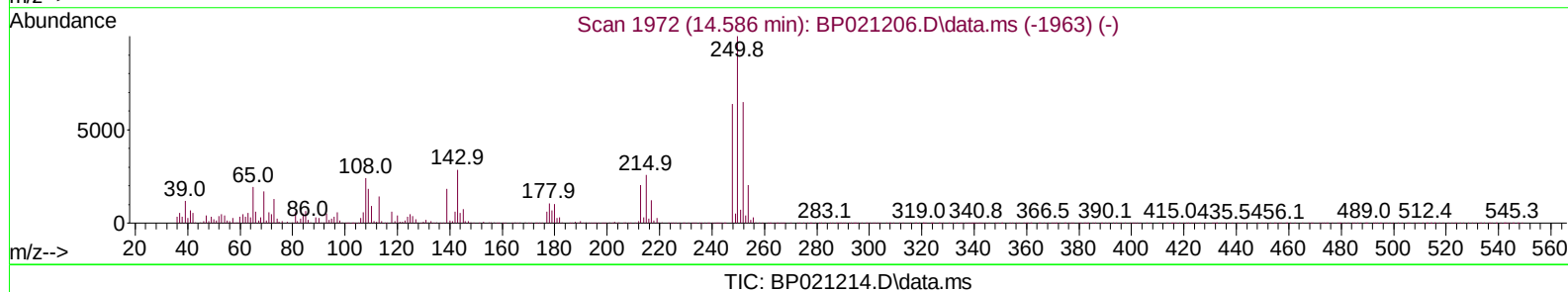
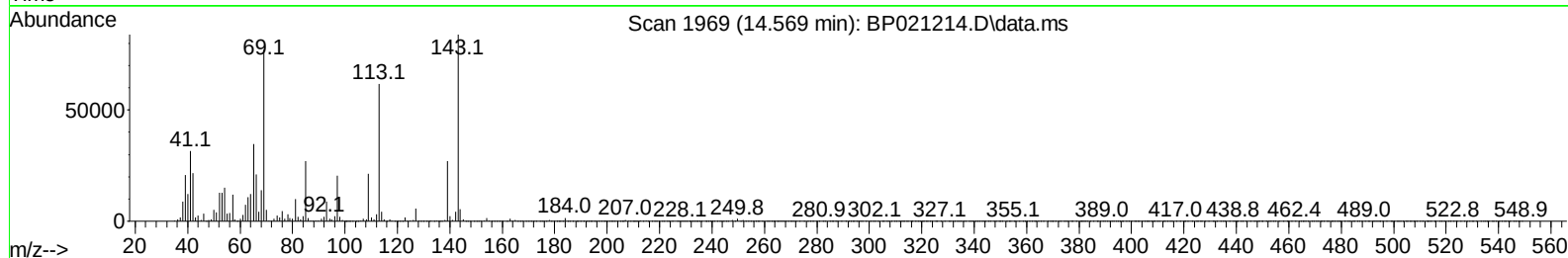
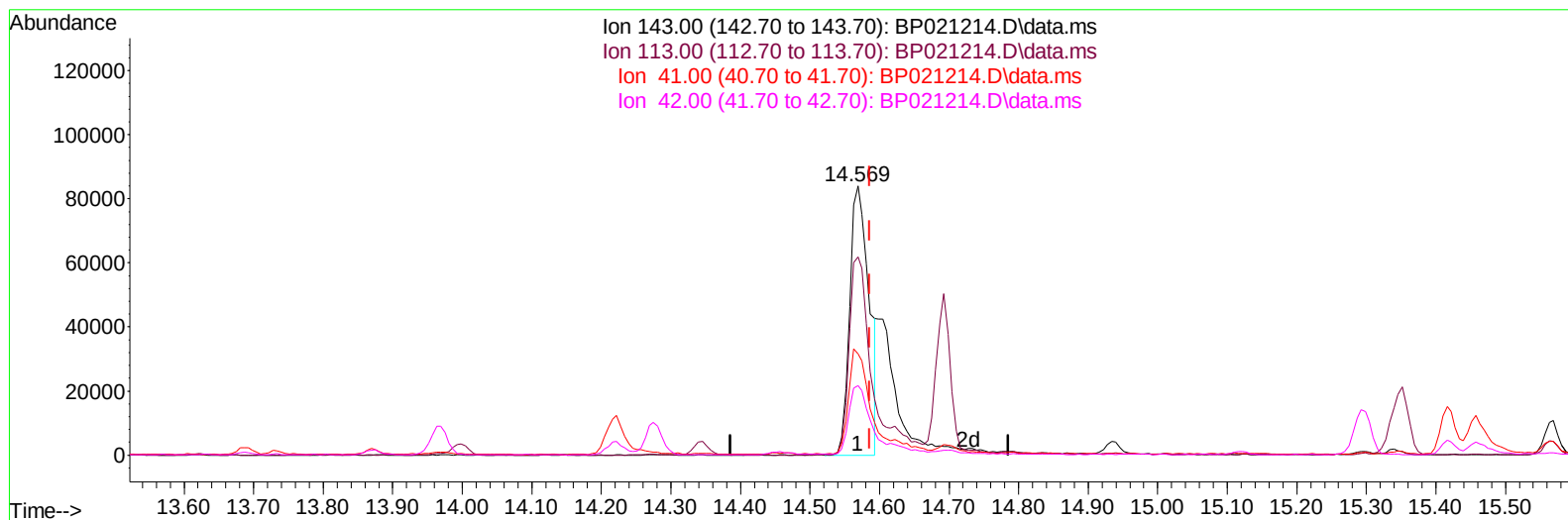
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Misc :
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Instrument :
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(54) 4-Nitrophenol-d4 (S)

14.569min (-0.017) 27.79 ng/ul

response 161365

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	73.58
41.00	37.90	37.73
42.00	27.10	25.85

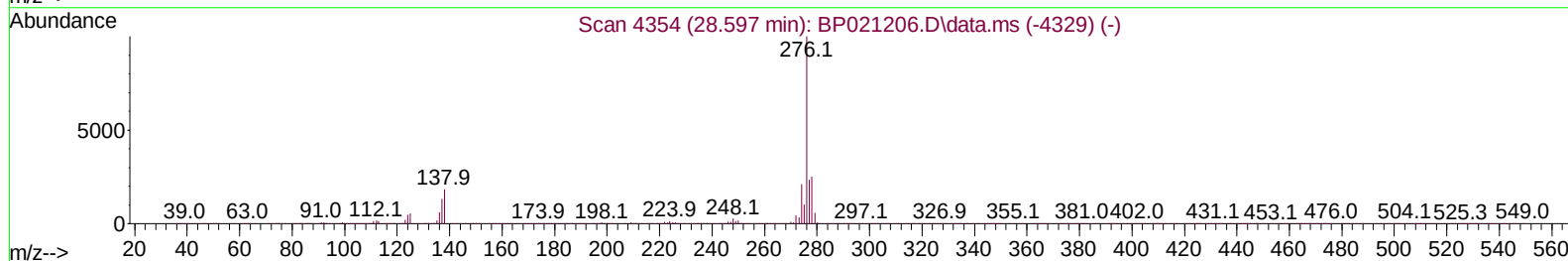
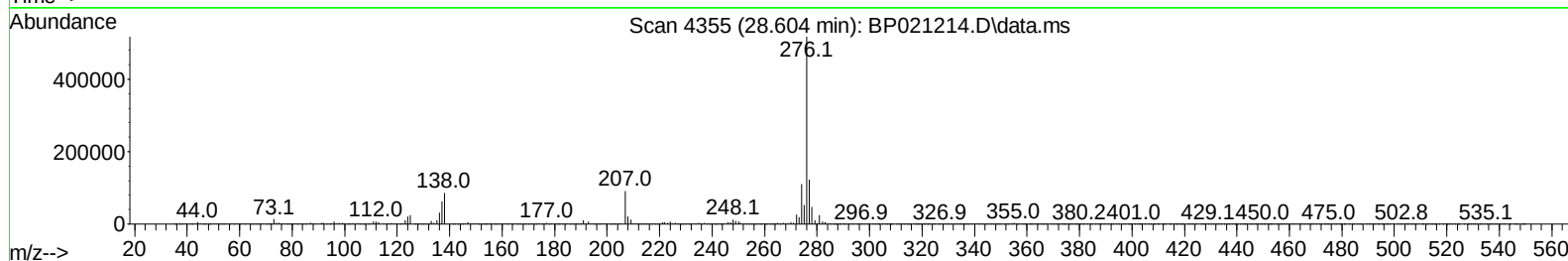
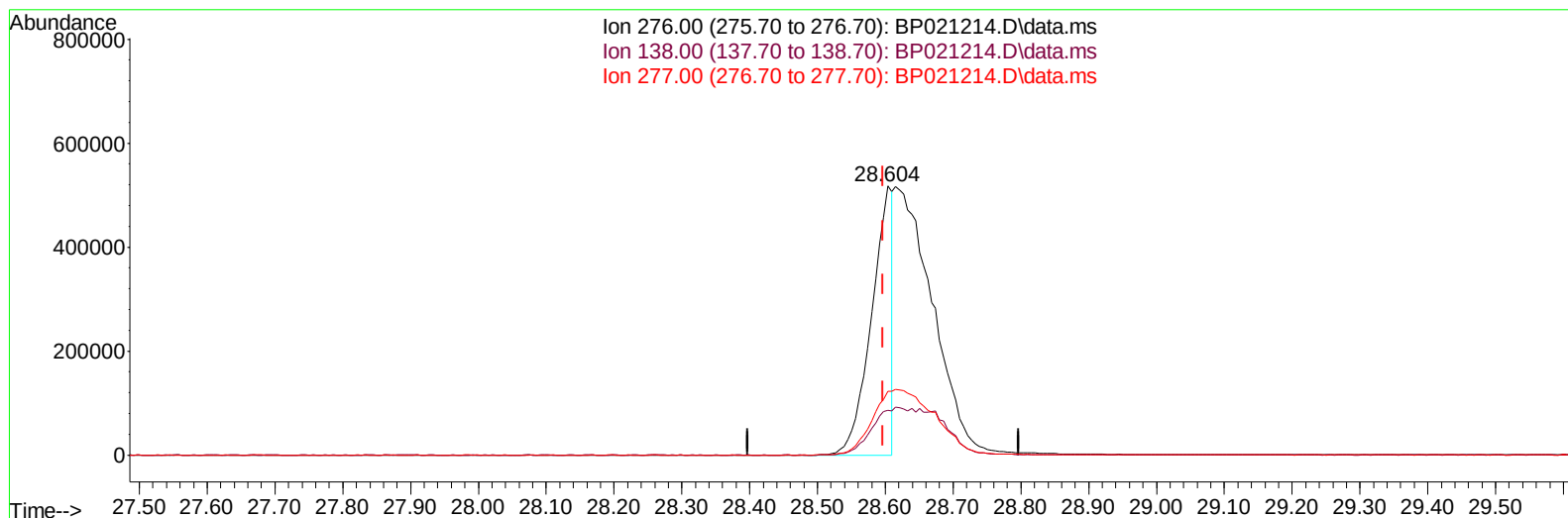
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021214.D
Acq On : 29 Jul 2024 17:50
Operator : MA/JU
Sample : PB162189BS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 07/30/2024
Supervised By :mohammad ahmed 09/04/2024



TIC: BP021214.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.604min (+ 0.006) 10.69 ng/ul

response 1116420

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	17.90	16.64
277.00	24.00	23.77
0.00	0.00	0.00

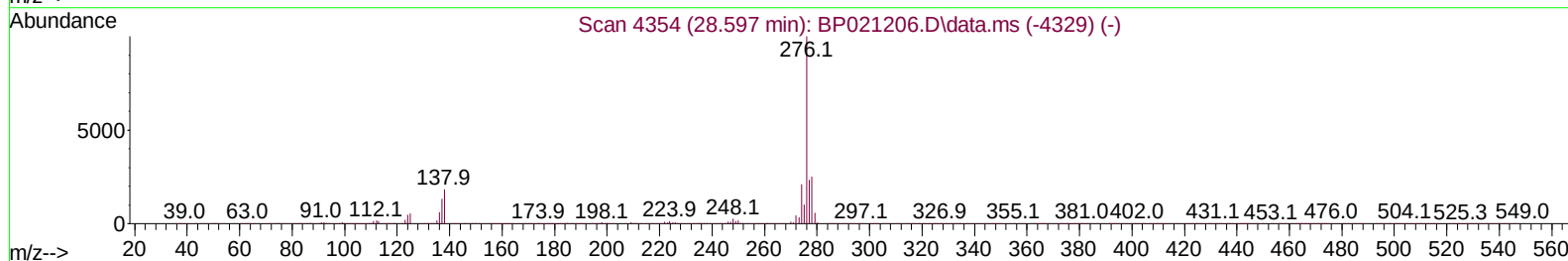
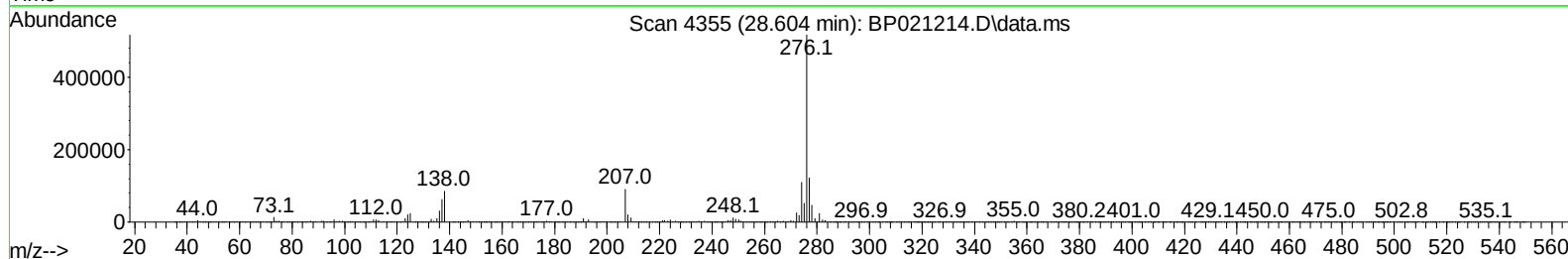
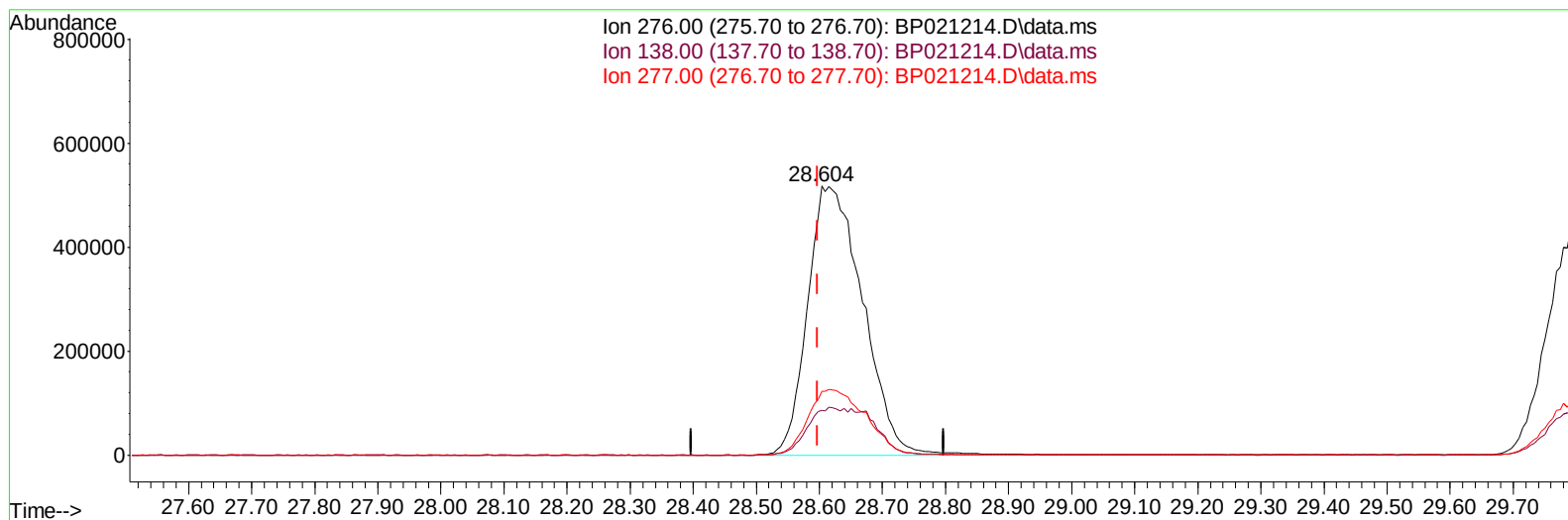
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021214.D
Acq On : 29 Jul 2024 17:50
Operator : MA/JU
Sample : PB162189BS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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TIC: BP021214.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.604min (+ 0.006) 29.92 ng/ul m

response 3124668

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	17.90	16.64
277.00	24.00	23.77
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Quant Time: Jul 29 19:11:30 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.634	152		220503	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136		863964	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164		561324	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188		1217338	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240		1220581	20.000	ng/ul	0.01
88) Perylene-d12	24.845	264		1413134	20.000	ng/ul	0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.170	96		30287	6.250	ng/uL	0.00
4) Pyridine-d5	3.587	84		364276	25.967	ng/ul	0.00
7) Phenol-d5	6.846	99		489399	27.224	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67		291358	26.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132		399232	26.954	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113		384213	25.733	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128		192325	28.660	ng/ul	0.00
24) 2-Nitrophenol-d4	9.499	143		222158	28.925	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165		398891	28.721	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131		476372	25.892	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166		1163704	28.518	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160		1314156	28.682	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143		239453m	41.244	ng/ul	-0.02
60) Fluorene-d10	15.298	176		1007433	30.094	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.457	200		220648	29.957	ng/ul	0.00
73) Anthracene-d10	17.198	188		1544215	28.561	ng/ul	0.00
81) Pyrene-d10	19.586	212		1964704	26.036	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.616	264		2117838	30.387	ng/ul	0.02
Target Compounds							
						Qvalue	
2) 1,4-Dioxane	3.205	88		54351	10.199	ng/uL	98
5) Pyridine	3.611	79		372915	25.773	ng/ul	99
6) Benzaldehyde	6.805	77		248224	27.389	ng/ul	97
8) Phenol	6.875	94		513511	28.229	ng/ul	99
10) Bis(2-Chloroethyl)ether	7.075	93		402166	26.762	ng/ul	99
12) 2-Chlorophenol	7.216	128		418402	28.185	ng/ul	98
13) 2-Methylphenol	8.093	108		375419	26.685	ng/ul	98
14) 2,2'-oxybis(1-Chloropr...	8.152	45		580226m	26.725	ng/ul	
16) Acetophenone	8.458	105		618909	26.858	ng/ul	95
17) N-Nitroso-di-n-propyla...	8.434	70		310412	25.708	ng/ul	98
18) 4-Methylphenol	8.428	108		409616	27.178	ng/ul	97
19) Hexachloroethane	8.681	117		180202	27.013	ng/ul	98
22) Nitrobenzene	8.834	77		471733	29.346	ng/ul	99
23) Isophorone	9.346	82		878576	27.492	ng/ul	100
25) 2-Nitrophenol	9.534	139		237815	29.688	ng/ul	98
26) 2,4-Dimethylphenol	9.599	107		348551	22.066	ng/ul	98
27) Bis(2-Chloroethoxy)met...	9.822	93		536205	27.604	ng/ul	99
29) 2,4-Dichlorophenol	10.075	162		398594	29.121	ng/ul	99
30) Naphthalene	10.446	128		1303372	28.680	ng/ul	99
32) 4-Chloroaniline	10.587	127		441812	25.147	ng/ul	99
33) Hexachlorobutadiene	10.699	225		289351	28.194	ng/ul	99
34) Caprolactam	11.387	113		133213m	31.594	ng/ul	
35) 4-Chloro-3-methylphenol	11.734	107		440936	29.637	ng/ul	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.063	142	880929	28.239	ng/ul	97
37) 1-Methylnaphthalene	12.281	142	881221	27.466	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.440	216	531860	29.176	ng/ul	97
40) Hexachlorocyclopentadiene	12.387	237	126372	13.252	ng/ul	97
41) 2,4,6-Trichlorophenol	12.704	196	293528	25.482	ng/ul	99
42) 2,4,5-Trichlorophenol	12.793	196	329777	27.059	ng/ul	99
43) 1,1'-Biphenyl	13.087	154	1175300	28.650	ng/ul	98
44) 2-Chloronaphthalene	13.134	162	944087	28.828	ng/ul	99
45) 2-Nitroaniline	13.369	65	291130	32.146	ng/ul	93
47) Dimethylphthalate	13.734	163	1200093	28.985	ng/ul	100
48) 2,6-Dinitrotoluene	13.875	165	253051	30.570	ng/ul	97
50) Acenaphthylene	13.998	152	1475887	28.450	ng/ul	100
51) 3-Nitroaniline	14.222	138	247455	34.353	ng/ul	97
52) Acenaphthene	14.345	153	972358	28.240	ng/ul	98
53) 2,4-Dinitrophenol	14.457	184	140332	32.282	ng/ul	95
55) 4-Nitrophenol	14.587	109	171631	36.342	ng/ul	96
56) Dibenzofuran	14.693	168	1380610	29.281	ng/ul	97
57) 2,4-Dinitrotoluene	14.704	165	373872	32.902	ng/ul	95
58) 2,3,4,6-Tetrachlorophenol	14.934	232	245492	23.523	ng/ul	98
59) Diethylphthalate	15.122	149	1217046	29.413	ng/ul	99
61) Fluorene	15.351	166	1146373	29.794	ng/ul	100
62) 4-Chlorophenyl-phenyle...	15.340	204	600137	28.744	ng/ul	98
63) 4-Nitroaniline	15.416	138	253896	42.562	ng/ul	94
66) 4,6-Dinitro-2-methylph...	15.475	198	237251	30.374	ng/ul	98
67) N-Nitrosodiphenylamine	15.569	169	988223	27.877	ng/ul	99
68) 4-Bromophenyl-phenylether	16.257	248	395392	27.711	ng/ul	98
69) Hexachlorobenzene	16.375	284	466715	28.527	ng/ul	97
70) Atrazine	16.563	200	23219	1.797	ng/ul	95
71) Pentachlorophenol	16.751	266	173323	19.506	ng/ul	97
72) Phenanthrene	17.139	178	1840479	29.144	ng/ul	100
74) Anthracene	17.239	178	1844901	28.657	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.051	216	516620	26.515	ng/uL	98
76) Pentachlorobenzene	14.604	250	514780	26.589	ng/uL	98
77) Carbazole	17.528	167	1721306	32.707	ng/ul	99
78) Di-n-butylphthalate	18.098	149	2178961	30.009	ng/ul	99
80) Fluoranthene	19.239	202	2300547	25.244	ng/ul	100
82) Pyrene	19.616	202	2436704	26.276	ng/ul	99
83) Butylbenzylphthalate	20.551	149	1041527	26.957	ng/ul	98
84) 3,3'-Dichlorobenzidine	21.445	252	862042	30.908	ng/ul	99
85) Benzo(a)anthracene	21.522	228	2443836	28.582	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.427	149	1472299	26.479	ng/ul	99
87) Chrysene	21.592	228	2363200	29.744	ng/ul	99
89) Di-n-octyl phthalate	22.686	149	2622713	28.095	ng/ul	100
90) Benzo(b)fluoranthene	23.798	252	2634248	30.718	ng/ul	99
91) Benzo(k)fluoranthene	23.869	252	2533515	29.603	ng/ul	99
93) Benzo(a)pyrene	24.698	252	2348931	28.986	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	28.604	276	3124668m	29.923	ng/ul	
95) Dibenzo(a,h)anthracene	28.674	278	2573851	29.765	ng/ul	99
96) Benzo(g,h,i)perylene	29.804	276	2444302	28.741	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Instrument :

BNA_P

ClientSampleId :

SLCS189

Manual IntegrationsAPPROVED

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