

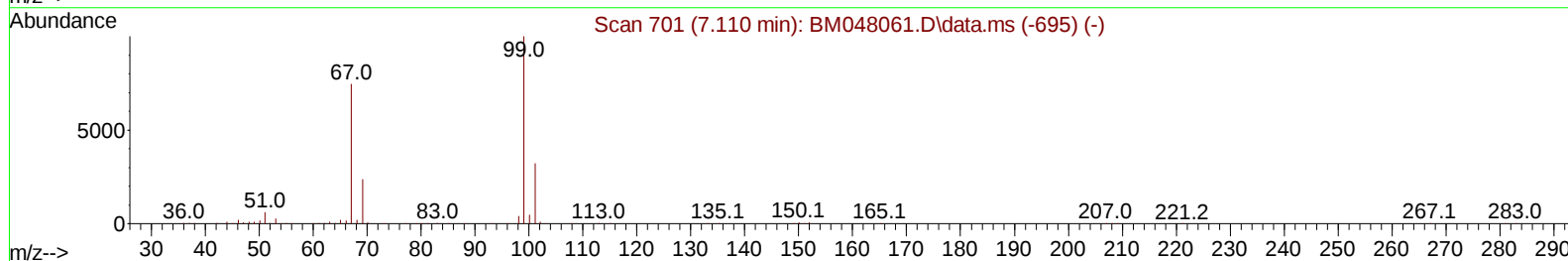
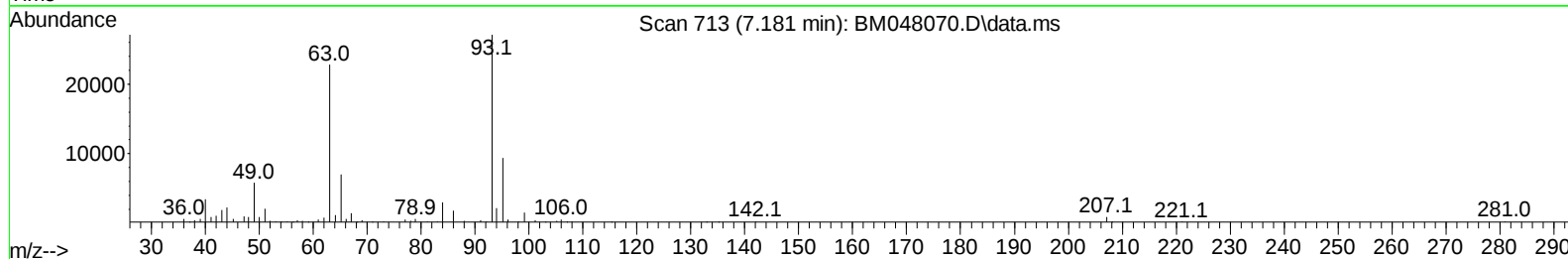
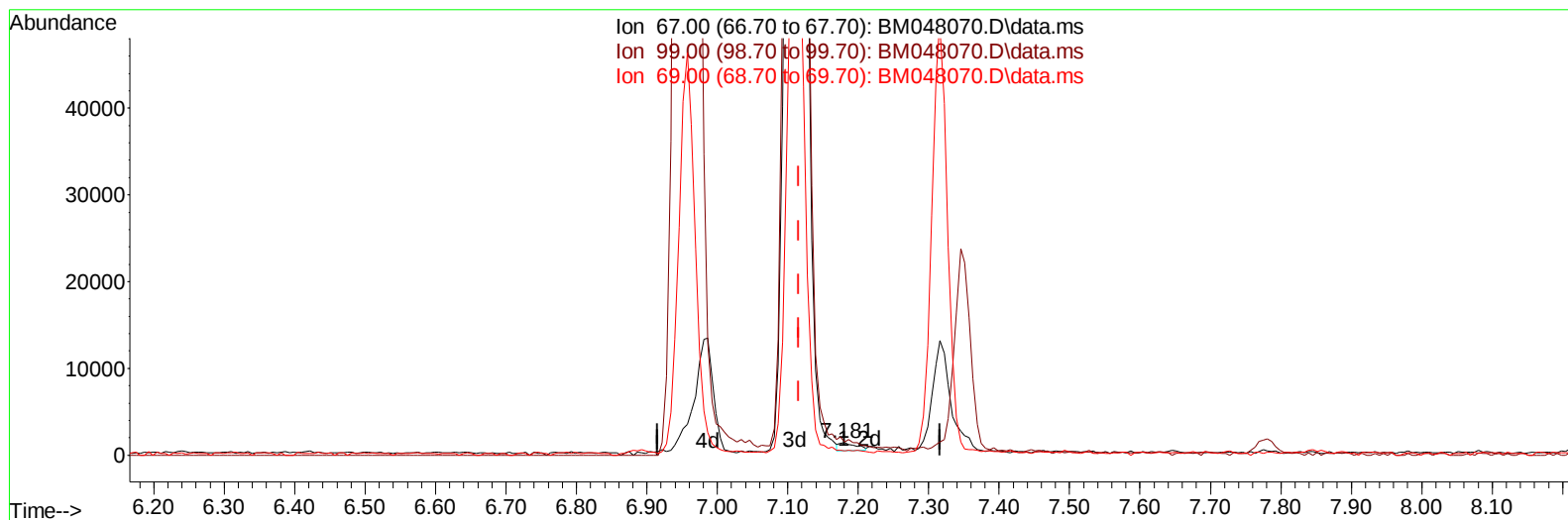
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048070.D
Acq On : 15 Oct 2024 21:05
Operator : RC/JU
Sample : P4350-14MSD
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOAK1MSD

Manual IntegrationsAPPROVED

Quant Time: Oct 15 23:03:38 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 10 05:10:14 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/16/2024
Supervised By :mohammad ahmed 10/17/2024



TIC: BM048070.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.181min (+ 0.065) 0.09 ng/u1

response 1569

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	106.00	105.56
69.00	32.00	32.68
0.00	0.00	0.00

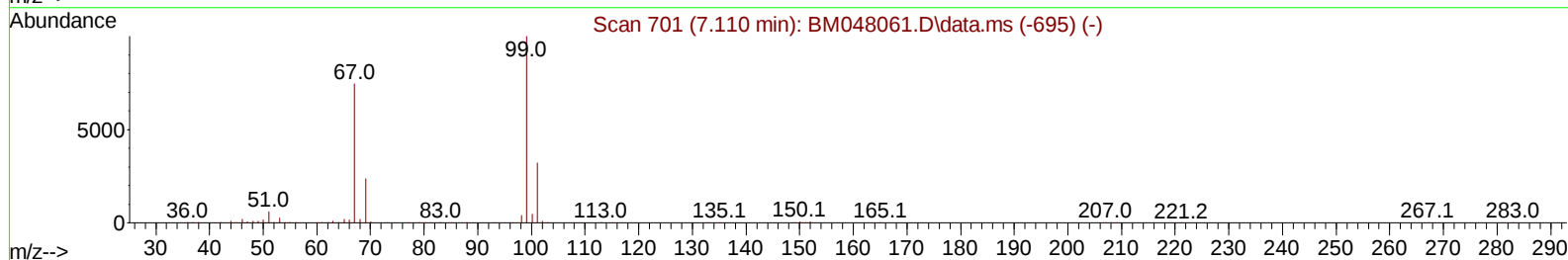
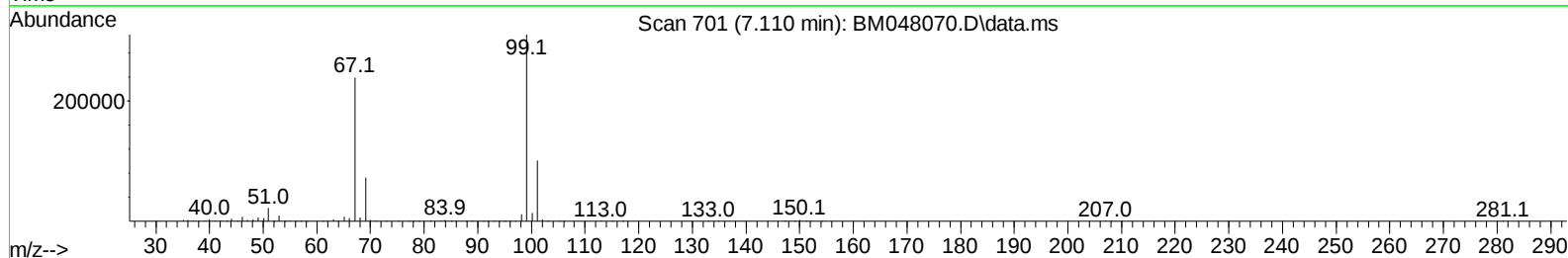
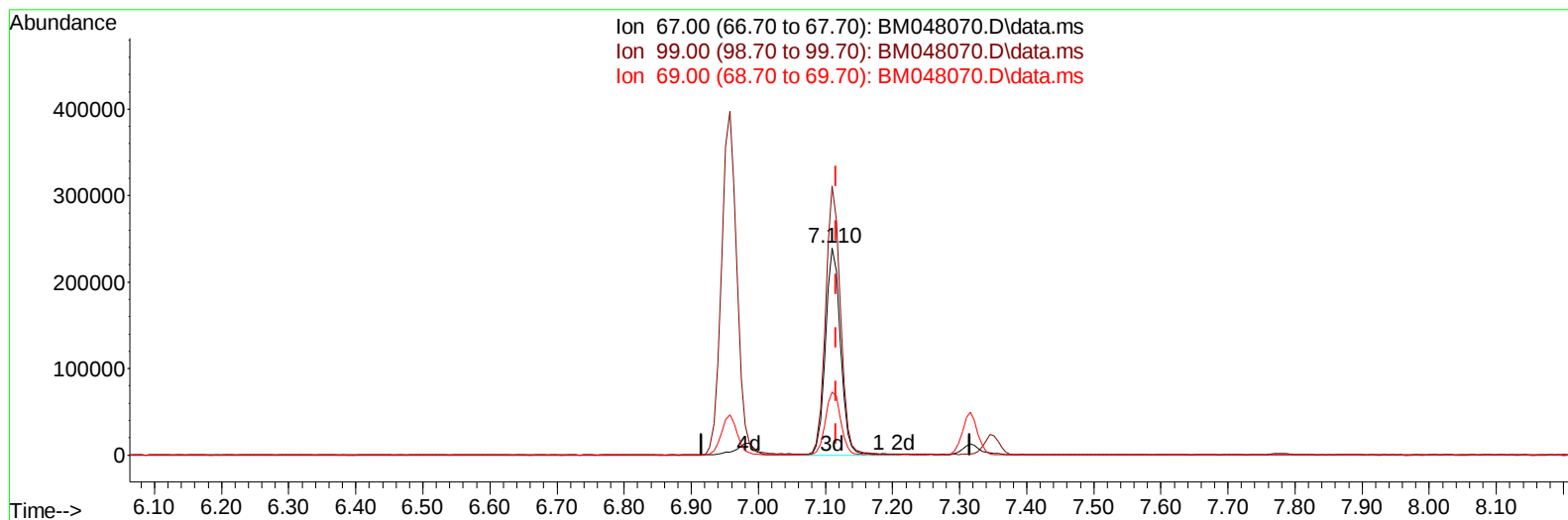
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(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.110min (-0.006) 20.46 ng/u1 m

response 373072

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	106.00	129.79#
69.00	32.00	30.62
0.00	0.00	0.00

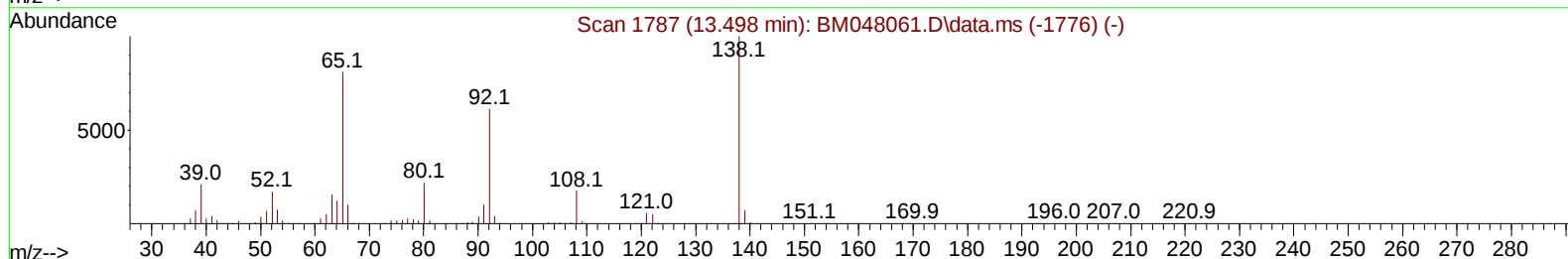
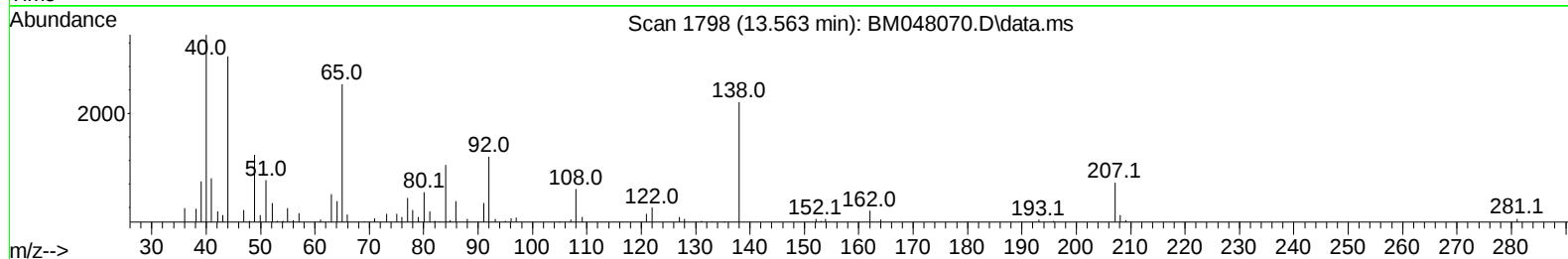
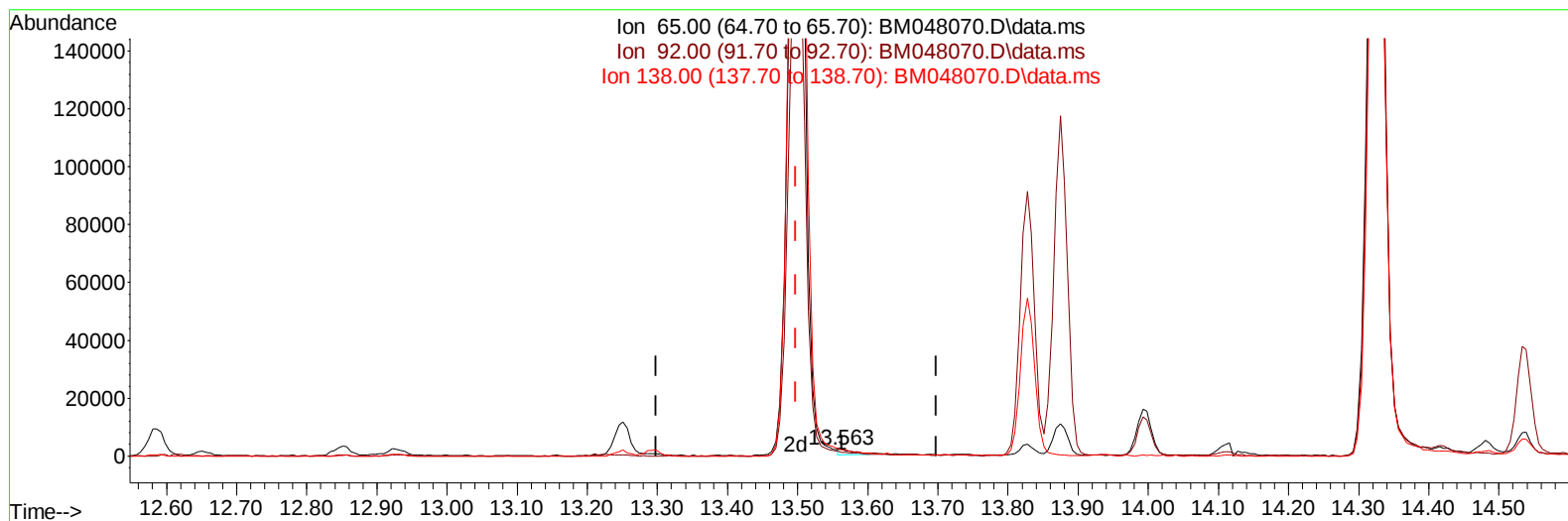
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TIC: BM048070.D\data.ms

(45) 2-Nitroaniline

13.563min (+ 0.065) 0.19 ng/ul

response 2642

Ion	Exp%	Act%
65.00	100.00	100.00
92.00	61.40	50.76
138.00	84.50	87.84
0.00	0.00	0.00

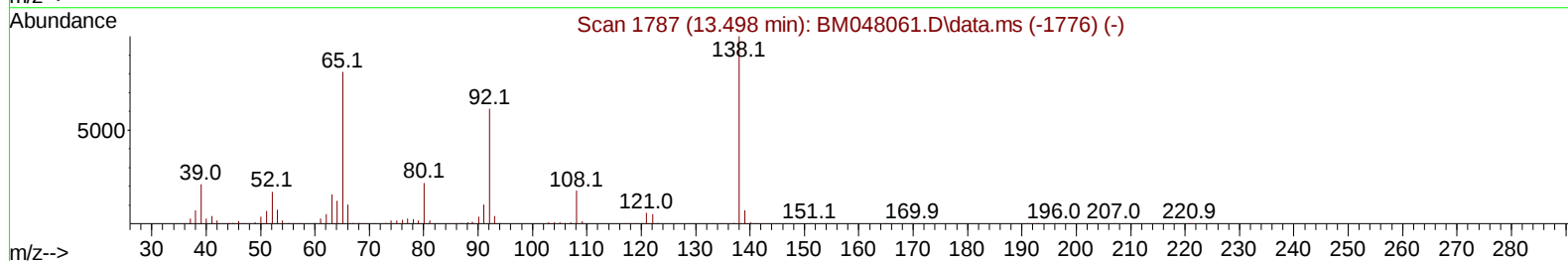
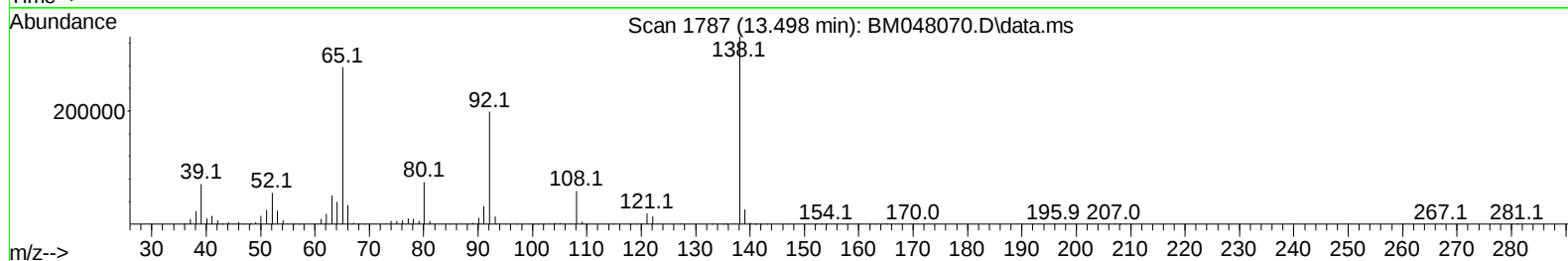
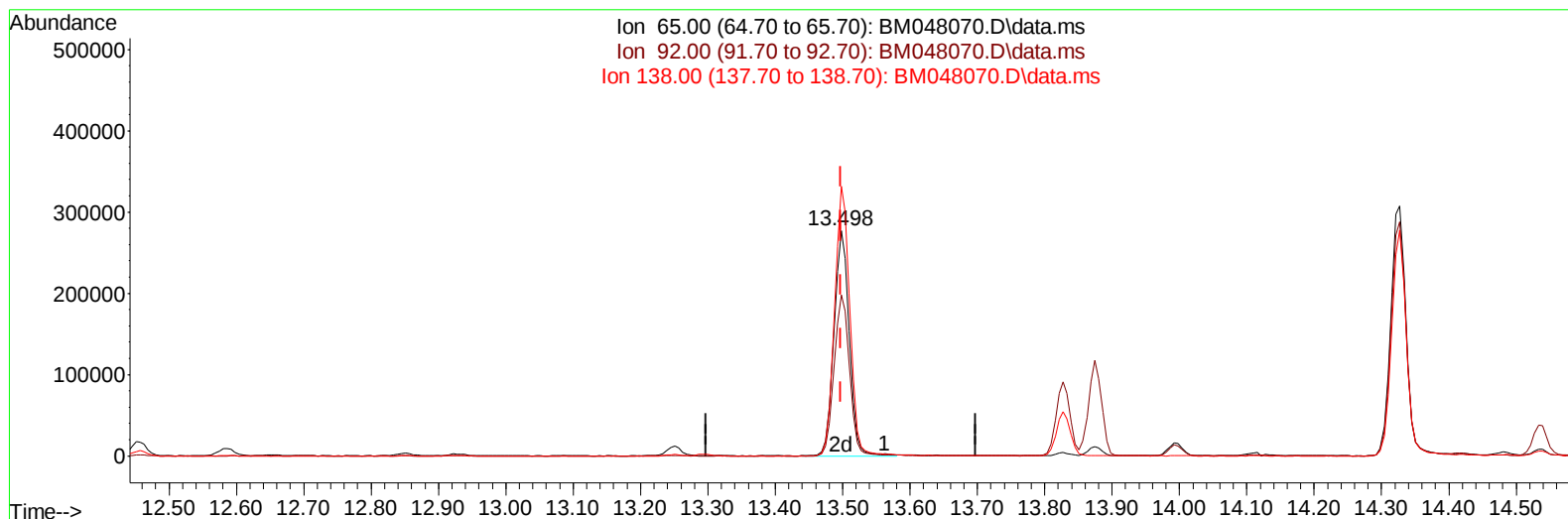
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Instrument :
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TIC: BM048070.D\data.ms

(45) 2-Nitroaniline

13.498min (+ 0.000) 32.15 ng/ul m

response 438548

Ion	Exp%	Act%
65.00	100.00	100.00
92.00	61.40	71.60
138.00	84.50	119.58#
0.00	0.00	0.00

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Instrument :
 BNA_M
 ClientSampleId :
 EOAK1MSD

Manual IntegrationsAPPROVED

Quant Time: Oct 15 23:14:06 2024
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 QLast Update : Thu Oct 10 05:10:14 2024
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Reviewed By :Yogesh Patel 10/16/2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	271691	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	1121286	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	693941	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.162	188	1369592	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240	1218750	20.000	ng/ul	0.00
88) Perylene-d12	24.380	264	1411183	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	31174	3.708	ng/uL	0.00
4) Pyridine-d5	3.652	84	458566	18.628	ng/ul	0.00
7) Phenol-d5	6.957	99	634362	24.692	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	373072m	20.464	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	531894	28.456	ng/ul	0.00
15) 4-Methylphenol-d8	8.492	113	521150	27.284	ng/ul	0.00
21) Nitrobenzene-d5	8.939	128	265570	29.747	ng/ul	0.00
24) 2-Nitrophenol-d4	9.657	143	306640	35.842	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	545537	31.920	ng/ul	0.00
31) 4-Chloroaniline-d4	10.710	131	698335	26.430	ng/ul	0.00
46) Dimethylphthalate-d6	13.827	166	1486911	29.554	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1728087	28.970	ng/ul	0.00
54) 4-Nitrophenol-d4	14.633	143	216645	21.628	ng/ul	0.01
60) Fluorene-d10	15.410	176	1254501	28.787	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	200881	25.064	ng/ul	0.00
73) Anthracene-d10	17.257	188	1882488	27.953	ng/ul	0.00
81) Pyrene-d10	19.551	212	2179139	31.504	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.180	264	2245549	30.669	ng/ul	0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.275	88	91184	9.930	ng/uL#	84
5) Pyridine	3.669	79	654453	25.253	ng/ul	86
6) Benzaldehyde	6.922	77	401936	28.768	ng/ul	87
8) Phenol	6.987	94	914782	33.464	ng/ul#	88
10) Bis(2-Chloroethyl)ether	7.204	93	673003	30.897	ng/ul	89
12) 2-Chlorophenol	7.351	128	717874	36.253	ng/ul#	90
13) 2-Methylphenol	8.228	108	651290	34.168	ng/ul	99
14) 2,2'-oxybis(1-Chloropr...	8.298	45	981402	23.934	ng/ul#	94
16) Acetophenone	8.604	105	1056891	32.337	ng/ul#	86
17) N-Nitroso-di-n-propyla...	8.592	70	506015	29.231	ng/ul#	85
18) 4-Methylphenol	8.557	108	724213	33.809	ng/ul	99
19) Hexachloroethane	8.857	117	294915	33.134	ng/ul	84
22) Nitrobenzene	8.981	77	768688	30.310	ng/ul#	87
23) Isophorone	9.510	82	1505880	33.193	ng/ul#	93
25) 2-Nitrophenol	9.686	139	421610	43.076	ng/ul#	82
26) 2,4-Dimethylphenol	9.757	107	610361	29.104	ng/ul	91
27) Bis(2-Chloroethoxy)met...	9.986	93	917823	33.077	ng/ul	98
29) 2,4-Dichlorophenol	10.228	162	693201	39.280	ng/ul	98
30) Naphthalene	10.622	128	2261533	35.064	ng/ul	99
32) 4-Chloroaniline	10.733	127	708702	27.240	ng/ul	100
33) Hexachlorobutadiene	10.910	225	431528	37.009	ng/ul	97
34) Caprolactam	11.522	113	229446	40.653	ng/ul#	69
35) 4-Chloro-3-methylphenol	11.869	107	702820	38.473	ng/ul	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.239	142	1521104	37.118	ng/ul	97
37) 1-Methylnaphthalene	12.457	142	1526153	36.604	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.610	216	813216	35.999	ng/ul#	96
40) Hexachlorocyclopentadiene	12.586	237	388891	28.463	ng/ul	99
41) 2,4,6-Trichlorophenol	12.851	196	535519	40.973	ng/ul	99
42) 2,4,5-Trichlorophenol	12.927	196	574564	40.200	ng/ul	96
43) 1,1'-Biphenyl	13.251	154	1964108	35.351	ng/ul#	97
44) 2-Chloronaphthalene	13.292	162	1544715	35.948	ng/ul	96
45) 2-Nitroaniline	13.498	65	438548m	32.149	ng/ul	
47) Dimethylphthalate	13.874	163	1892082	37.134	ng/ul	98
48) 2,6-Dinitrotoluene	13.998	165	401797	44.902	ng/ul#	77
50) Acenaphthylene	14.139	152	2434699	36.191	ng/ul	99
51) 3-Nitroaniline	14.327	138	427045	39.431	ng/ul	81
52) Acenaphthene	14.480	153	1629190	34.259	ng/ul	97
53) 2,4-Dinitrophenol	14.539	184	241258	43.633	ng/ul#	68
55) 4-Nitrophenol	14.645	109	251229	29.872	ng/ul	84
56) Dibenzofuran	14.816	168	2261703	35.332	ng/ul	97
57) 2,4-Dinitrotoluene	14.786	165	569162	42.349	ng/ul#	75
58) 2,3,4,6-Tetrachlorophenol	15.051	232	449826	39.927	ng/ul#	93
59) Diethylphthalate	15.239	149	1894268	37.157	ng/ul	98
61) Fluorene	15.468	166	1757327	32.906	ng/ul	100
62) 4-Chlorophenyl-phenyle...	15.463	204	875933	34.711	ng/ul	94
63) 4-Nitroaniline	15.492	138	443383	42.154	ng/ul	87
66) 4,6-Dinitro-2-methylph...	15.551	198	350924	39.515	ng/ul#	83
67) N-Nitrosodiphenylamine	15.674	169	1522181	35.666	ng/ul	98
68) 4-Bromophenyl-phenylether	16.351	248	549819	38.946	ng/ul	97
69) Hexachlorobenzene	16.474	284	642002	39.077	ng/ul	91
70) Atrazine	16.627	200	367857	25.893	ng/ul	99
71) Pentachlorophenol	16.815	266	387271	37.071	ng/ul	100
72) Phenanthrene	17.204	178	2780416	34.618	ng/ul	99
74) Anthracene	17.292	178	2784231	33.988	ng/ul	100
75) 1,2,3,4-Tetrachloroben...	13.216	216	788040	34.751	ng/uL	99
76) Pentachlorobenzene	14.739	250	761915	34.422	ng/uL	98
77) Carbazole	17.562	167	2620530	35.292	ng/ul	98
78) Di-n-butylphthalate	18.121	149	3384289	40.939	ng/ul#	98
80) Fluoranthene	19.215	202	3202025	39.305	ng/ul	99
82) Pyrene	19.580	202	3408324	37.590	ng/ul	95
83) Butylbenzylphthalate	20.468	149	1536699	47.939	ng/ul#	84
84) 3,3'-Dichlorobenzidine	21.292	252	1008695	39.169	ng/ul	98
85) Benzo(a)anthracene	21.368	228	3283632	38.480	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.286	149	2177777	45.251	ng/ul	97
87) Chrysene	21.433	228	3067189	36.870	ng/ul	99
89) Di-n-octyl phthalate	22.415	149	3973196	43.651	ng/ul	100
90) Benzo(b)fluoranthene	23.438	252	3333580	37.879	ng/ul	98
91) Benzo(k)fluoranthene	23.497	252	3294765	36.614	ng/ul	99
93) Benzo(a)pyrene	24.244	252	3020146	36.135	ng/ul	98
94) Indeno(1,2,3-cd)pyrene	27.768	276	3966450	37.032	ng/ul	98
95) Dibenzo(a,h)anthracene	27.821	278	3213300	36.607	ng/ul	98
96) Benzo(g,h,i)perylene	28.826	276	3089843	36.819	ng/ul	97

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

Instrument :

BNA_M

ClientSampleId :

EOAK1MSD

Manual IntegrationsAPPROVED

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