

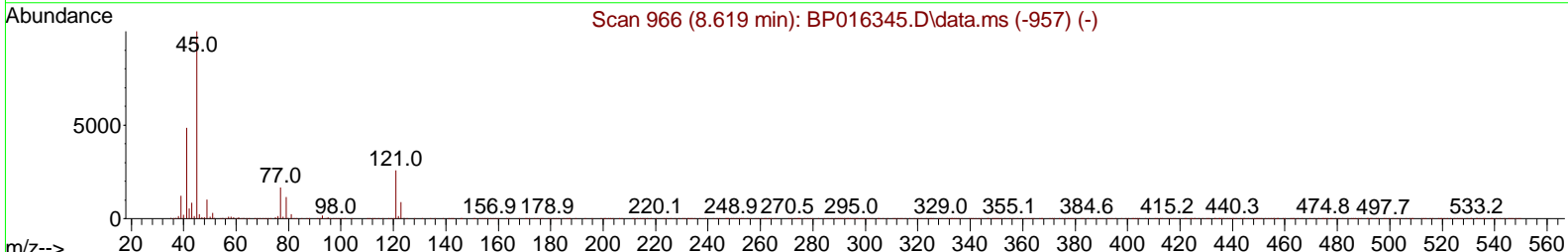
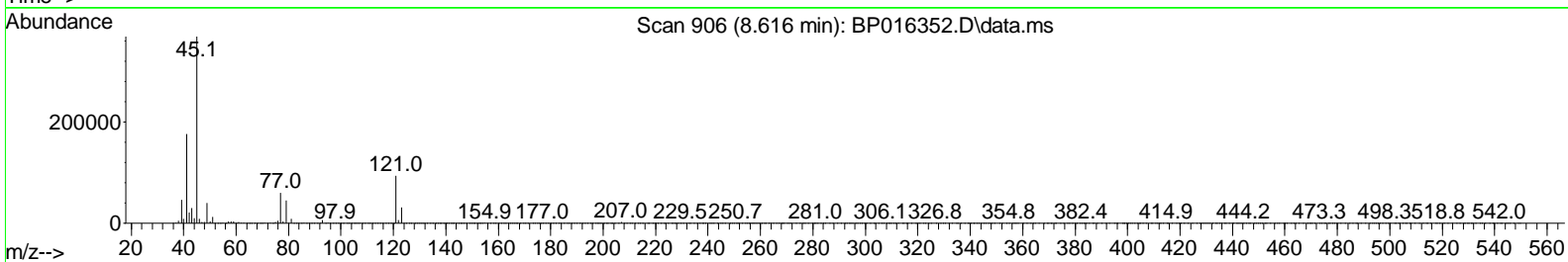
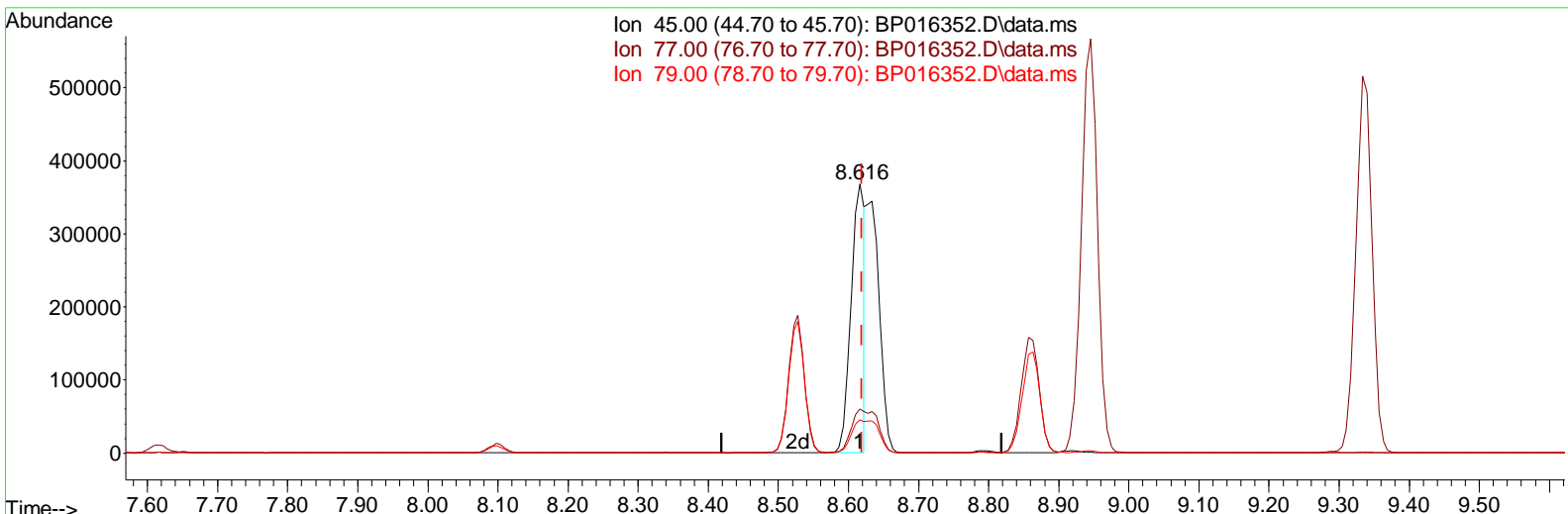
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072523\
 Data File : BP016352.D
 Acq On : 25 Jul 2023 18:49
 Operator : MA/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
LabSampleID :
 SSTDCCC020

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 07/26/2023
 Supervised By :mohammad ahmed 07/26/2023

Quant Time: Jul 26 00:29:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 25 23:56:16 2023
 Response via : Initial Calibration



TIC: BP016352.D\data.ms

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

8.616min (-0.003) 10.72 ng/ul

response 498126

Ion	Exp%	Act%
45.00	100.00	100.00
77.00	17.10	16.37
79.00	12.00	12.33
0.00	0.00	0.00

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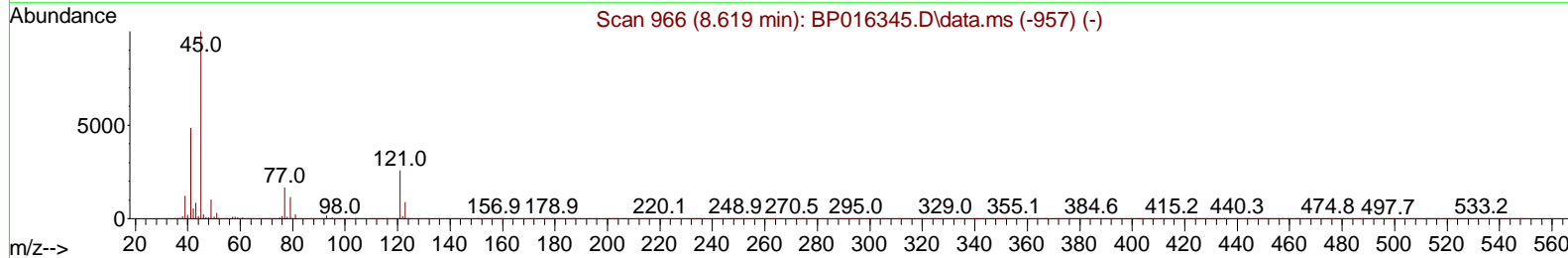
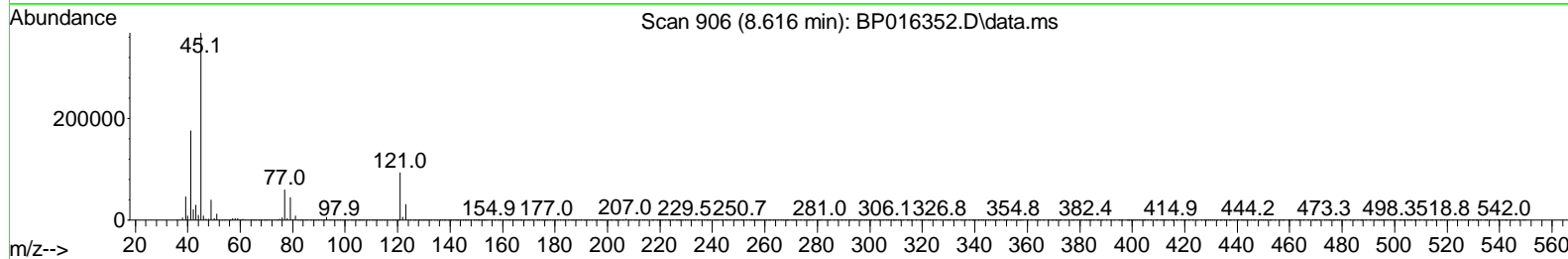
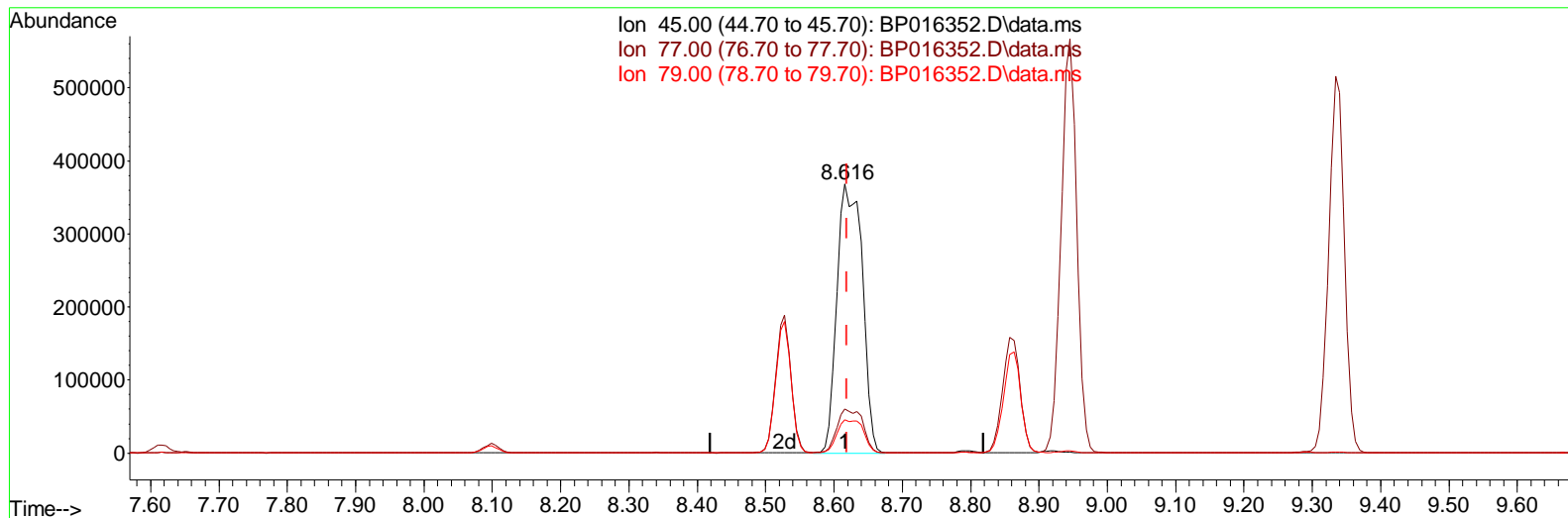
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(14) 2,2'-oxybis(1-Chloropropane)

8.616min (-0.003) 20.37 ng/ul m

response 946169

Ion	Exp%	Act%
45.00	100.00	100.00
77.00	17.10	16.37
79.00	12.00	12.33
0.00	0.00	0.00

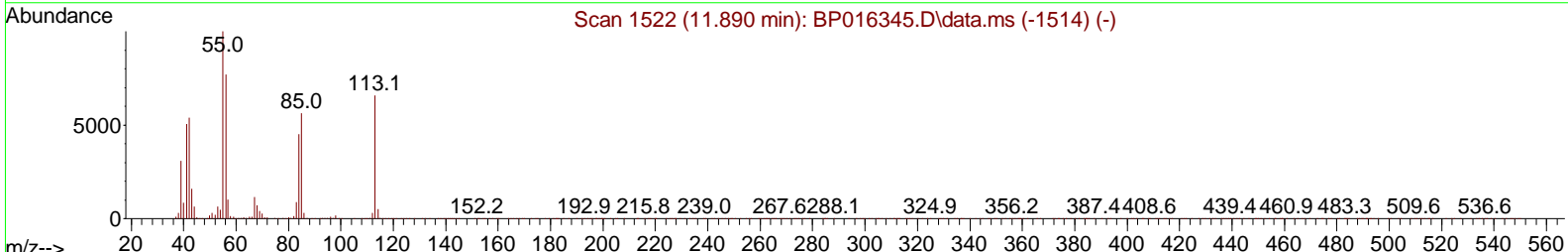
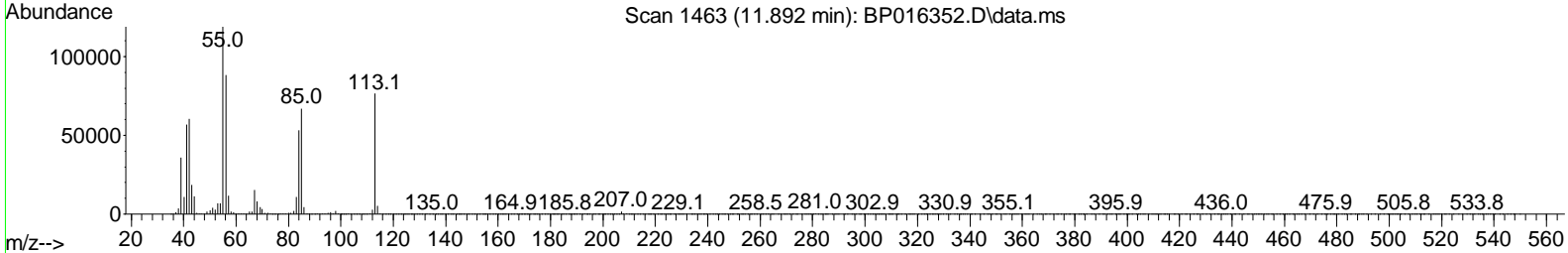
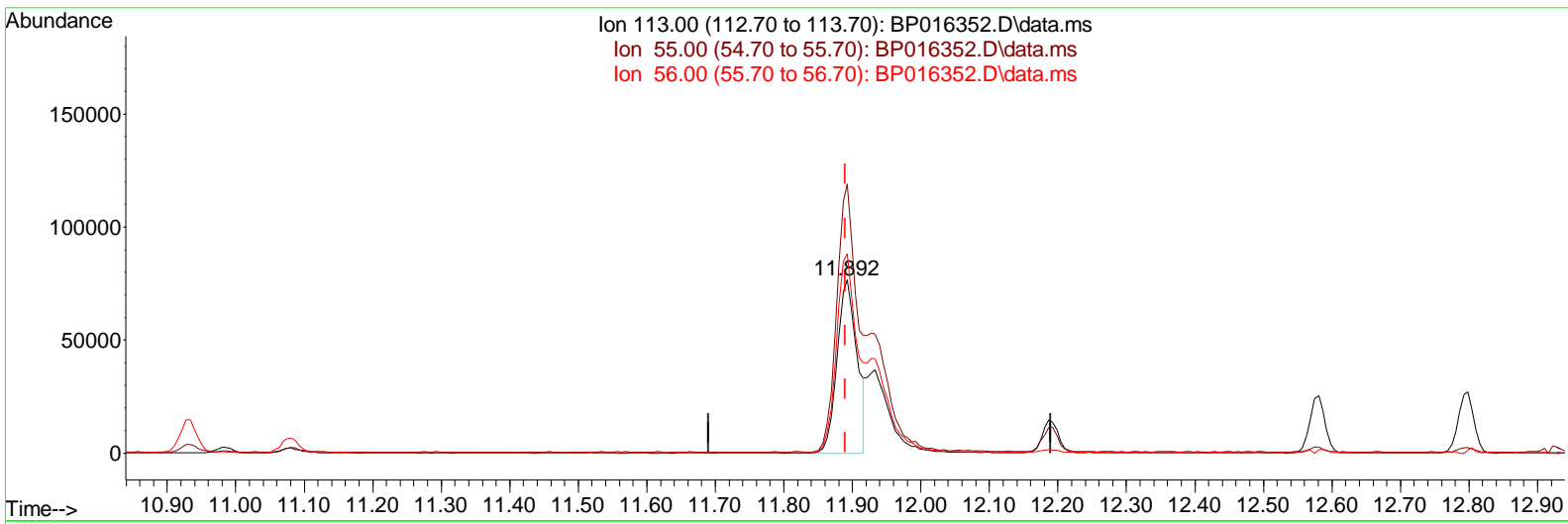
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TIC: BP016352.D\data.ms

(34) Caprolactam

11.892min (+ 0.003) 13.32 ng/ul

response 155674

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	152.80	155.33
56.00	117.30	115.09
0.00	0.00	0.00

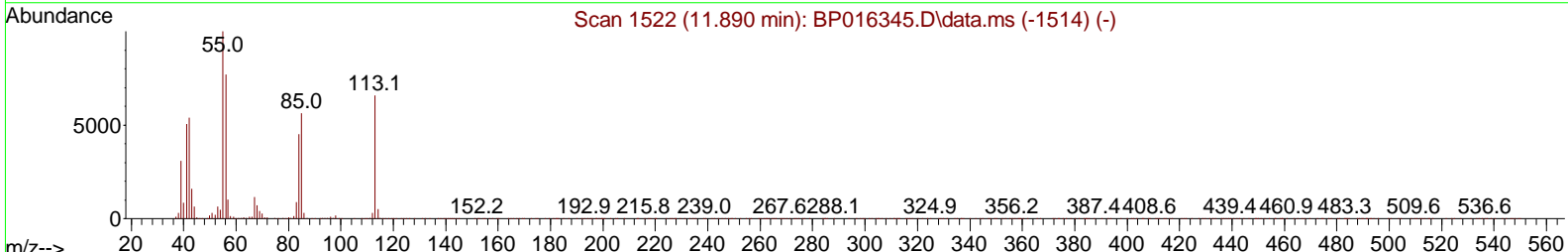
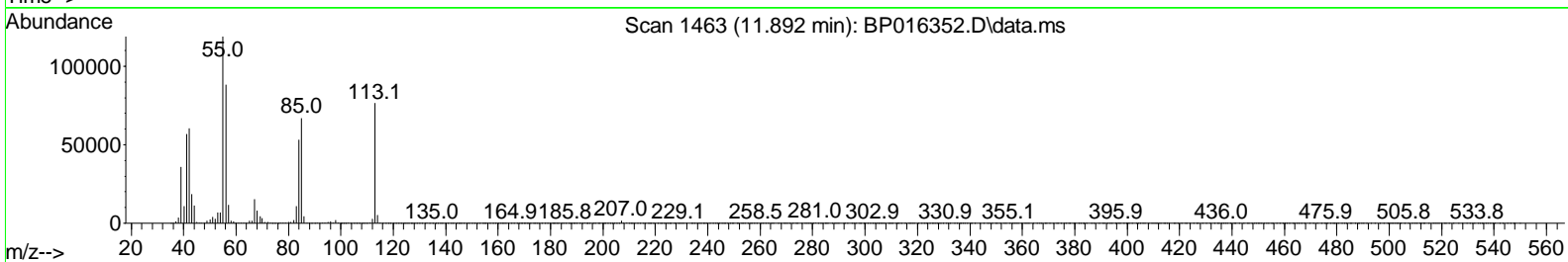
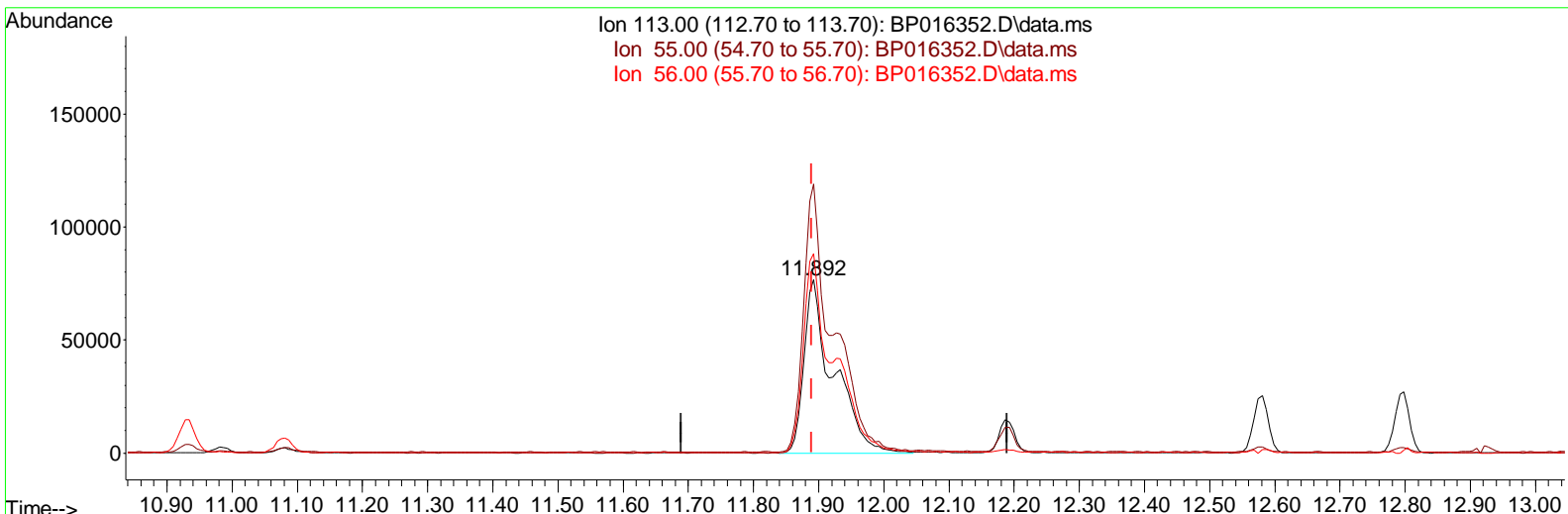
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TIC: BP016352.D\data.ms

(34) Caprolactam

11.892min (+ 0.003) 20.55 ng/ul m

response 240062

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	152.80	155.33
56.00	117.30	115.09
0.00	0.00	0.00

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LabSampleID :
 SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/26/2023
 Supervised By :mohammad ahmed 07/26/2023

Quant Time: Jul 26 00:36:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 25 23:56:16 2023
 Response via : Initial Calibration

Compound	R.T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.098	152	538586	20.000	ng/ul	0.00
20) Naphthalene-d8	10.934	136	2304811	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.739	164	1414049	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188	3033460	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	2549971	20.000	ng/ul	0.00
88) Perylene-d12	24.204	264	2481078	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.428	96	110397	8.082	ng/uL	0.00
4) Pyridine-d5	3.858	84	770626	19.885	ng/ul	0.00
7) Phenol-d5	7.228	99	891774	19.786	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.434	67	545867	20.173	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	703685	20.013	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	723718	19.967	ng/ul	0.00
21) Nitrobenzene-d5	9.293	128	325248	22.917	ng/ul	0.00
24) 2-Nitrophenol-d4	10.010	143	286583	24.484	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	677839	20.892	ng/ul	0.00
31) 4-Chloroaniline-d4	11.081	131	1029593	20.398	ng/ul	0.00
46) Dimethylphthalate-d6	14.151	166	2070403	20.006	ng/ul	0.00
49) Acenaphthylene-d8	14.439	160	2449617	20.035	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	357284	20.604	ng/ul	0.00
60) Fluorene-d10	15.733	176	1736899	19.728	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.845	200	190189	19.035	ng/ul	0.00
73) Anthracene-d10	17.604	188	2642717	19.667	ng/ul	0.00
81) Pyrene-d10	19.833	212	2926715	20.214	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.033	264	2351314	19.643	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.469	88	115978	8.007	ng/uL	97
5) Pyridine	3.875	79	792291	20.094	ng/ul	99
6) Benzaldehyde	7.246	77	566128	23.024	ng/ul	99
8) Phenol	7.251	94	934093	19.930	ng/ul	99
10) Bis(2-Chloroethyl)ether	7.528	93	764854	19.920	ng/ul	100
12) 2-Chlorophenol	7.646	128	725687	20.073	ng/ul	100
13) 2-Methylphenol	8.528	108	711555	19.799	ng/ul	99
14) 2,2'-oxybis(1-Chloropr...	8.616	45	946169m	20.365	ng/ul	
16) Acetophenone	8.946	105	1181726	20.659	ng/ul	100
17) N-Nitrosodipropylamine	8.916	70	601837	20.205	ng/ul	98
18) 4-Methylphenol	8.857	108	762330	19.853	ng/ul	98
19) Hexachloroethane	9.181	117	302650	20.010	ng/ul	97
22) Nitrobenzene	9.334	77	828823	21.814	ng/ul	99
23) Isophorone	9.851	82	1725666	20.164	ng/ul	99
25) 2-Nitrophenol	10.040	139	349426	23.924	ng/ul	98
26) 2,4-Dimethylphenol	10.075	107	841291	20.372	ng/ul	99
27) Bis(2-Chloroethoxy)meth...	10.340	93	1034624	20.051	ng/ul	100
29) 2,4-Dichlorophenol	10.557	162	677495	20.656	ng/ul	99
30) Naphthalene	10.987	128	2374592	19.892	ng/ul	100
32) 4-Chloroaniline	11.104	127	1003509	20.280	ng/ul	99
33) Hexachlorobutadiene	11.222	225	426780	19.798	ng/ul	99
34) Caprolactam	11.892	113	240062m	20.547	ng/ul	
35) 4-Chloro-3-methylphenol	12.192	107	751062	20.687	ng/ul	99

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Instrument :
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LabSampled :
 SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/26/2023
 Supervised By :mohammad ahmed 07/26/2023

Quant Time: Jul 26 00: 36: 04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
 Quant Title : SVOA CALI BRATI ON
 QLast Update : Tue Jul 25 23: 56: 16 2023
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
36) 2-Methyl naphthal ene	12. 581	142	1619364	19. 905	ng/ul	100
37) 1-Methyl naphthal ene	12. 798	142	1648869	20. 272	ng/ul	99
39) 1, 2, 4, 5-Tetrachl oroben. . .	12. 922	216	828505	19. 832	ng/ul	99
40) Hexachl orocycl opentadi ene	12. 881	237	499208	16. 739	ng/ul	97
41) 2, 4, 6-Tri chl orophenol	13. 169	196	519069	21. 089	ng/ul	98
42) 2, 4, 5-Tri chl orophenol	13. 234	196	564581	21. 433	ng/ul	99
43) 1, 1' -Bi phenyl	13. 581	154	2099461	19. 663	ng/ul	99
44) 2-Chl oronaphthal ene	13. 628	162	1644459	19. 688	ng/ul	99
45) 2-Ni troani li ne	13. 845	65	426163	24. 384	ng/ul	96
47) Di methyl phthal ate	14. 198	163	2090192	20. 028	ng/ul	99
48) 2, 6-Di ni trotol uene	14. 333	165	367955	24. 466	ng/ul	95
50) Acenaphthyl ene	14. 469	152	2764876	19. 850	ng/ul	100
51) 3-Ni troani li ne	14. 669	138	389428	22. 133	ng/ul #	98
52) Acenaphthene	14. 804	153	1763717	19. 637	ng/ul	100
53) 2, 4-Di ni trophenol	14. 863	184	103429	16. 080	ng/ul	95
55) 4-Ni trophenol	14. 939	109	289508	20. 712	ng/ul	99
56) Di benzofuran	15. 139	168	2413212	19. 627	ng/ul	100
57) 2, 4-Di ni trotol uene	15. 110	165	522802	24. 382	ng/ul	98
58) 2, 3, 4, 6-Tetrachl orophenol	15. 345	232	450579	21. 442	ng/ul	95
59) Di ethyl phthal ate	15. 557	149	2101284	19. 990	ng/ul	99
61) Fl uorene	15. 792	166	1968501	19. 710	ng/ul	98
62) 4-Chl orophenyl -phenyl e. . .	15. 786	204	974416	20. 014	ng/ul	99
63) 4-Ni troani li ne	15. 833	138	365029	22. 625	ng/ul	92
66) 4, 6-Di ni tro-2-methyl ph. . .	15. 863	198	239842	20. 079	ng/ul	98
67) N-Ni trosodi phenyl ami ne	16. 004	169	1685878	19. 821	ng/ul	99
68) 4-Bromophenyl -phenyl ether	16. 686	248	574866	19. 983	ng/ul	98
69) Hexachl orobenzene	16. 763	284	681084	19. 743	ng/ul	100
70) Atrazi ne	16. 957	200	625230	20. 148	ng/ul	100
71) Pentachl orophenol	17. 122	266	382904	19. 697	ng/ul	99
72) Phenanthrene	17. 551	178	3073010	19. 468	ng/ul	100
74) Anthracene	17. 639	178	3151681	19. 601	ng/ul	100
75) 1, 2, 3, 4-Tetrachl oroben. . .	13. 534	216	869538	19. 858	ng/uL	99
76) Pentachl orobenzene	15. 033	250	769040	19. 875	ng/uL	99
77) Carbazol e	17. 916	167	2811833	20. 316	ng/ul	100
78) Di -n-butyl phthal ate	18. 439	149	3550867	20. 146	ng/ul	100
80) Fl uoranthene	19. 504	202	3469393	19. 989	ng/ul	99
82) Pyrene	19. 857	202	3593545	20. 041	ng/ul	100
83) Butyl benzyl phthal ate	20. 727	149	1436872	21. 166	ng/ul	99
84) 3, 3' -Di chl orobenzi di ne	21. 515	252	1038787	18. 535	ng/ul	99
85) Benzo(a)anthracene	21. 580	228	3361031	19. 460	ng/ul	100
86) Bi s(2-ethyl hexyl)phtha. . .	21. 480	149	2104249	20. 321	ng/ul	100
87) Chrysene	21. 639	228	3118405	19. 497	ng/ul	100
89) Di -n-octyl phthal ate	22. 486	149	3311593	21. 234	ng/ul	100
90) Benzo(b)fl uoranthene	23. 392	252	3012355	20. 274	ng/ul	100
91) Benzo(k)fl uoranthene	23. 445	252	2971233	20. 094	ng/ul	99
93) Benzo(a)pyrene	24. 086	252	2765759	19. 765	ng/ul	99
94) I ndeno(1, 2, 3-cd)pyrene	26. 986	276	2807386	17. 374	ng/ul	98
95) Di benzo(a, h)anthracene	27. 021	278	2371929	17. 748	ng/ul	99
96) Benzo(g, h, i)peryl ene	27. 845	276	2148164	16. 792	ng/ul	99

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

Instrument :

BNA_P

LabSampleId :

SSTDCCC020

Manual IntegrationsAPPROVED

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