

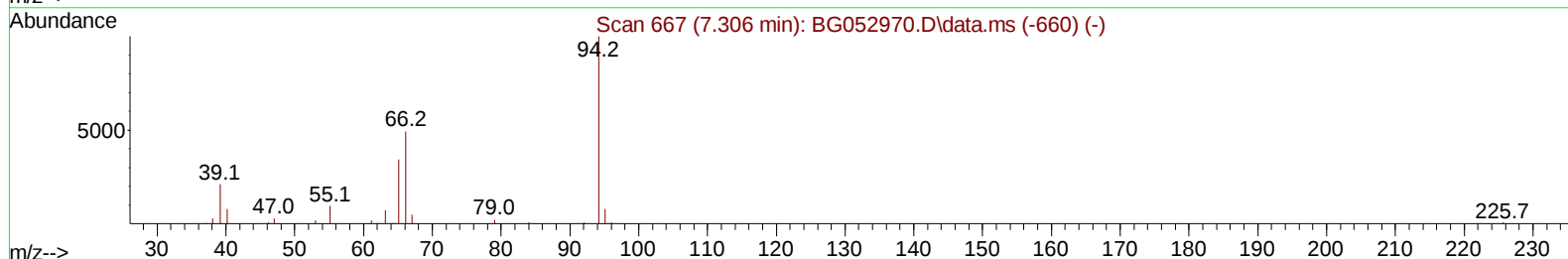
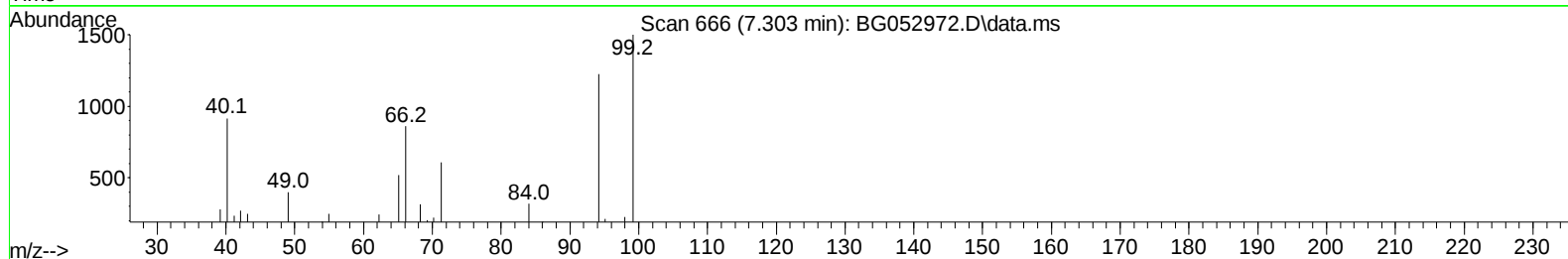
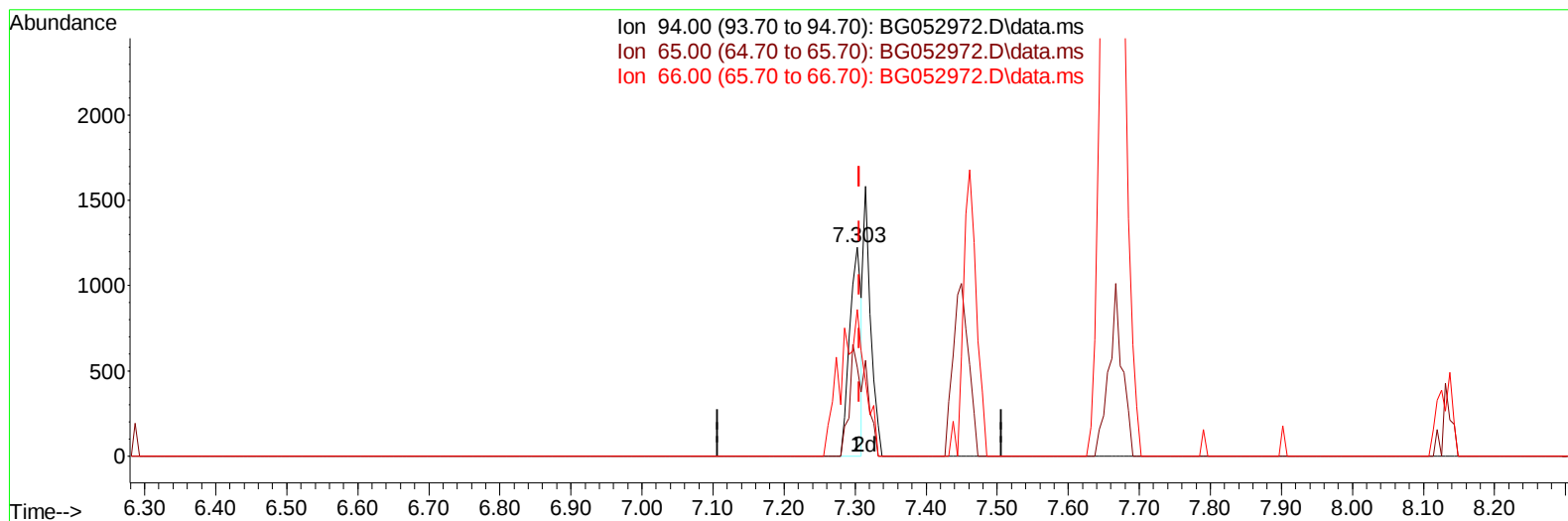
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052972.D
Acq On : 31 Mar 2022 9:56
Operator : CG/JU
Sample : N2160-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAYZ8

Manual IntegrationsAPPROVED

Quant Time: Apr 01 01:01:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/01/2022
Supervised By :Yogesh Patel 04/06/2022



TIC: BG052972.D\data.ms

(8) Phenol

7.303min (-0.004) 0.63 ng/u1

response 1434

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	32.40	42.50#
66.00	44.30	70.23#
0.00	0.00	0.00

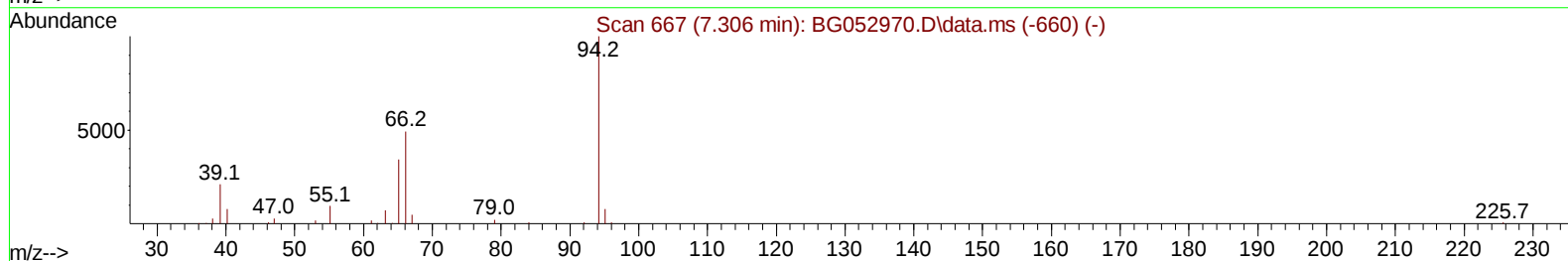
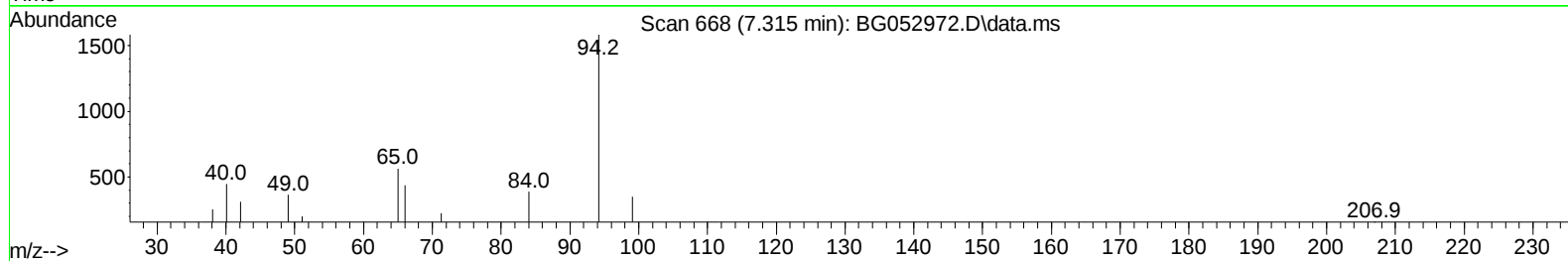
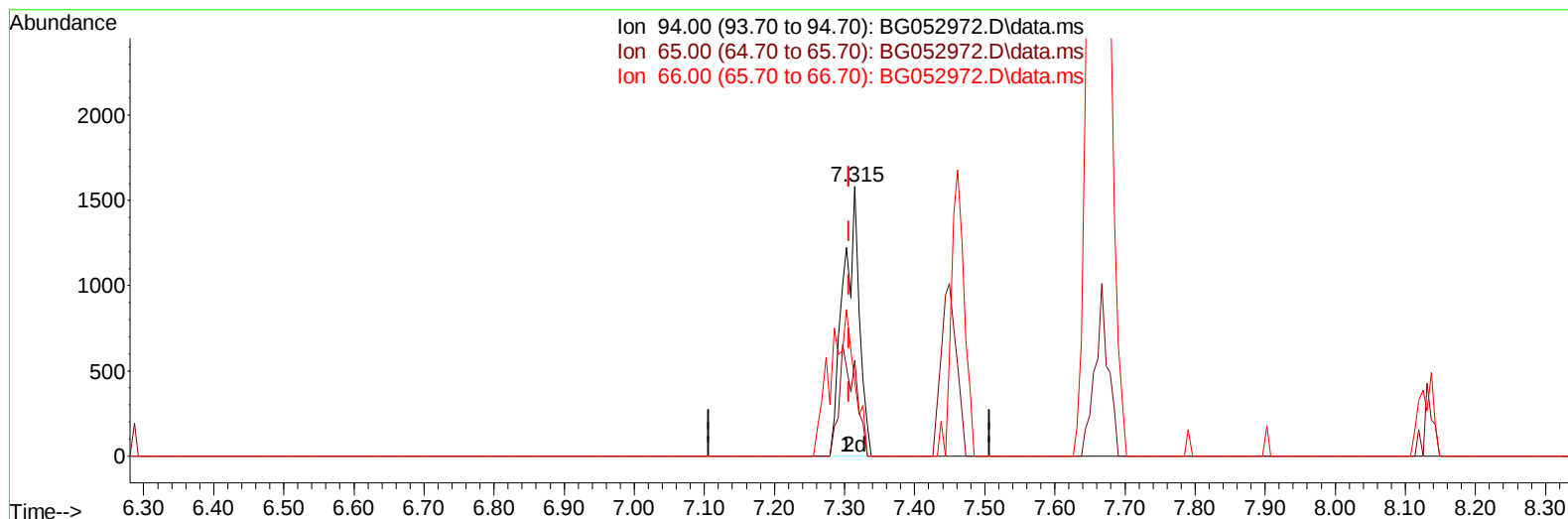
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TIC: BG052972.D\data.ms

(8) Phenol

7.315min (+ 0.008) 1.09 ng/ul m

response 2507

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	32.40	35.59
66.00	44.30	27.75#
0.00	0.00	0.00

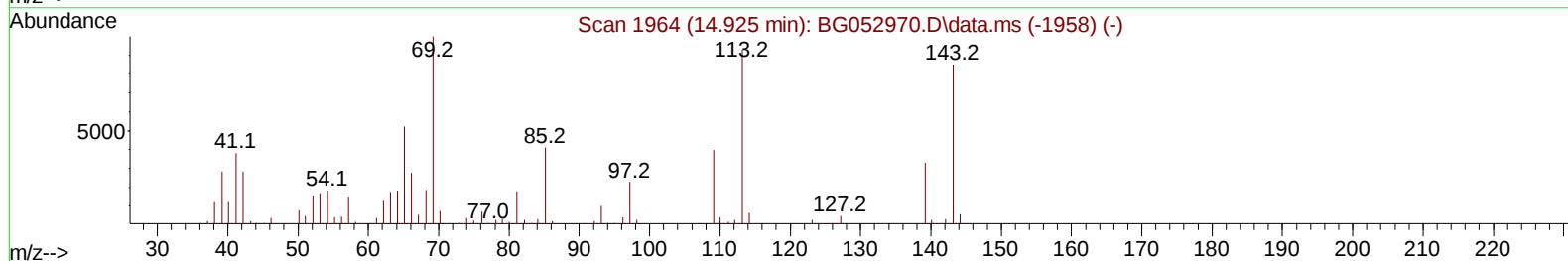
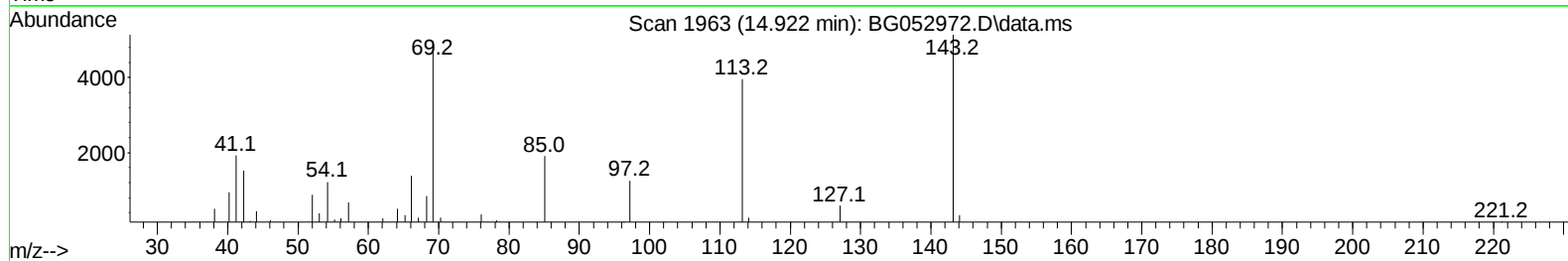
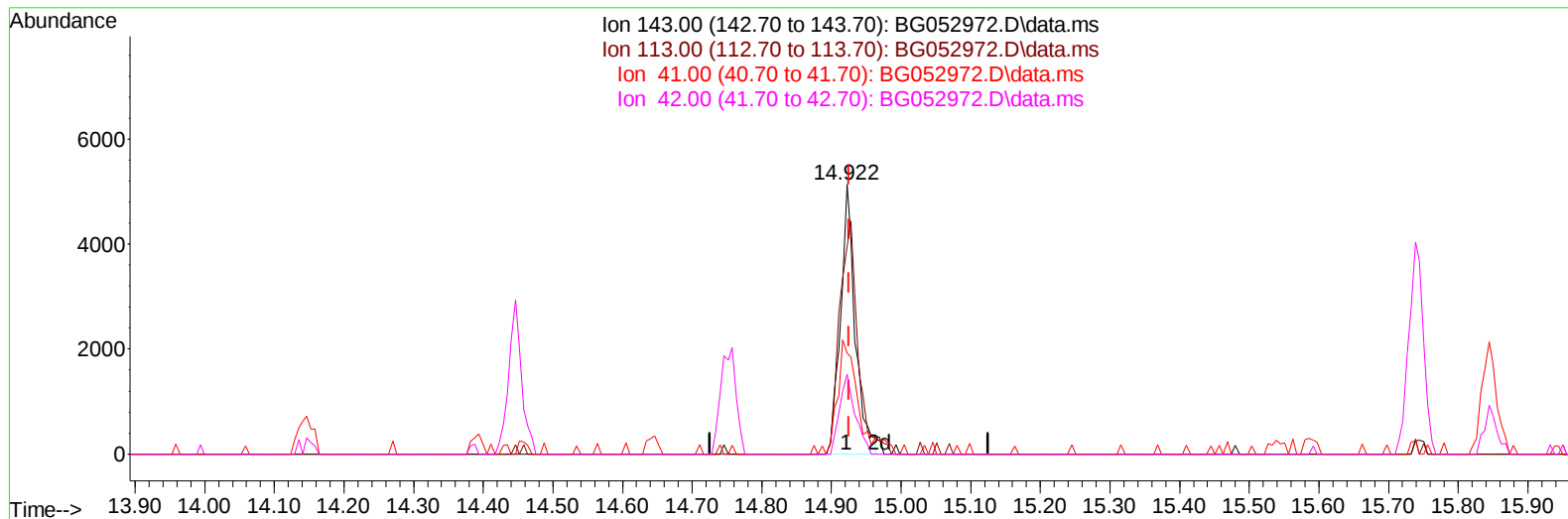
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Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BG052972.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (-0.004) 6.76 ng/u1

response 7615

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	76.94#
41.00	47.20	37.35#
42.00	29.40	29.66

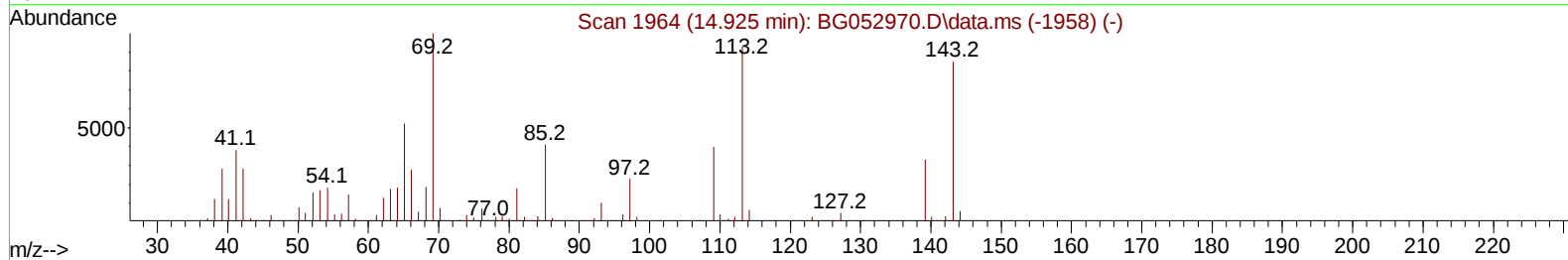
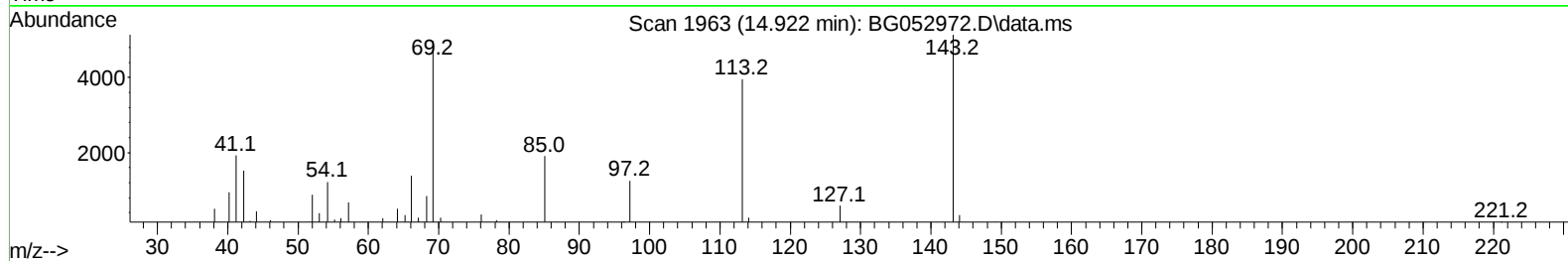
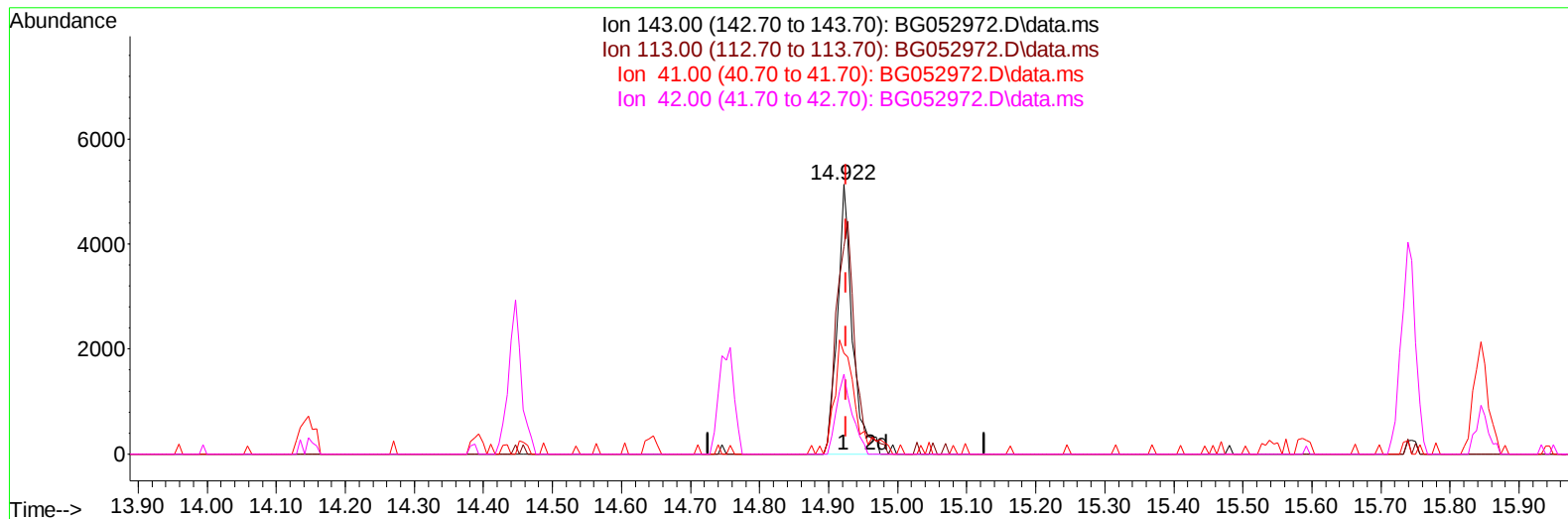
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 Sample : N2160-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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TIC: BG052972.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (-0.004) 6.86 ng/ul m

response 7730

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	76.94#
41.00	47.20	37.35#
42.00	29.40	29.66

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.137	152	26562	20.000	ng/ul	0.00
20) Naphthalene-d8	10.945	136	112873	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.752	164	87254	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.495	188	211453	20.000	ng/ul	0.00
79) Chrysene-d12	21.777	240	231629	20.000	ng/ul	# 0.00
88) Perylene-d12	25.085	264	232297	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.549	96	2606	3.466	ng/uL	0.00
4) Pyridine-d5	3.978	84	8636	4.600	ng/ul	0.01
7) Phenol-d5	7.279	99	16886	7.098	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.450	67	50915	35.041	ng/ul	0.00
11) 2-Chlorophenol-d4	7.661	132	39957	24.735	ng/ul	0.00
15) 4-Methylphenol-d8	8.824	113	30299	17.581	ng/ul	0.00
21) Nitrobenzene-d5	9.288	128	29347	36.119	ng/ul	0.00
24) 2-Nitrophenol-d4	10.017	143	31621	36.898	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	51927	27.852	ng/ul	0.00
31) 4-Chloroaniline-d4	11.068	131	58154	23.601	ng/ul	0.00
46) Dimethylphthalate-d6	14.147	166	234136	34.929	ng/ul	0.00
49) Acenaphthylene-d8	14.446	160	251577	30.886	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	7730m	6.859	ng/ul	0.00
60) Fluorene-d10	15.739	176	202533	34.411	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.844	200	38693	35.957	ng/ul	0.00
73) Anthracene-d10	17.595	188	346303	35.170	ng/ul	0.00
81) Pyrene-d10	19.874	212	433172	33.429	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.861	264	430899	35.900	ng/ul	0.00
Target Compounds						
8) Phenol	7.315	94	2507m	1.093	ng/ul	Qvalue

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

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