

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060120\
 Data File : BM026233.D
 Acq On : 02 Jun 2020 18:29
 Operator : JU/CG
 Sample : L2829-06DL 25X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5CD0DL

Quant Time: Jun 02 19:01:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 02 03:00:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	138809	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	548279	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.12	164	315350	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.87	188	690606	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	754065	20.00	ng/ul	0.00
86) Perylene-d12	23.18	264	794047	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.68	99	2737	0.22	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.83	67	9296	1.07	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	8634	0.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	5728	0.60	ng/ul	0.00
19) Nitrobenzene-d5	8.62	128	4631	1.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.33	143	3946	0.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.87	165	7724	0.87	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.54	166	31400	1.24	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	34159	1.14	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.12	176	28527	1.30	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	16.96	188	48198	1.42	ng/ul	0.00
79) Pyrene-d10	19.27	212	53004	1.37	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.05	264	46635	1.10	ng/ul	0.00

Target Compounds

					Ovalue	
40) 2,4,5-Trichlorophenol	12.62	196	8144	1.122	ng/ul#	92
56) 2,3,4,6-Tetrachlorophenol	14.76	232	10543	1.728	ng/ul#	96
69) Pentachlorophenol	16.53	266	220541	44.284	ng/ul	97

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

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