

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002025.D
 Acq On : 03 Mar 2020 17:19
 Operator : CG/JU
 Sample : L1617-08
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDS2

Quant Time: Mar 04 07:12:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 03 05:31:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	295882	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1230587	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	778669	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1695604	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1697235	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1869065	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	32381	4.70	ng/uL	0.00
5) Phenol-d5	6.82	99	578083	25.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	332712	27.84	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	514125	26.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	472669	25.48	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	238914	25.27	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	269014	25.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	514084	25.09	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	686287	30.56	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1501049	25.47	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1843230	26.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	202215	17.95	ng/ul	0.00
58) Fluorene-d10	15.27	176	1331247	26.11	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	149047	15.94	ng/ul	0.00
71) Anthracene-d10	17.13	188	1998391	26.22	ng/ul	0.00
79) Pyrene-d10	19.37	212	2289740	25.77	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2564243	26.82	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.84	94	45997	1.963	ng/ul 96
45) Dimethylphthalate	13.73	163	209283	3.440	ng/ul 100

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

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