

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002016.D
 Acq On : 03 Mar 2020 10:11
 Operator : CG/JU
 Sample : L1617-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDR7

Quant Time: Mar 04 05:12:51 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	284673	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1210742	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	765256	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1652222	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1623680	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1816535	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	44247	6.68	ng/uL	0.00
5) Phenol-d5	6.82	99	763192	35.15	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	439317	38.21	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	680646	36.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	623719	34.95	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	316661	34.05	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	349740	33.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	665109	33.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	570406	25.81	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1958354	33.81	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2379168	34.35	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	262350	23.70	ng/ul	0.00
58) Fluorene-d10	15.27	176	1753450	34.99	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	204708	22.47	ng/ul	0.00
71) Anthracene-d10	17.13	188	2592204	34.91	ng/ul	0.00
79) Pyrene-d10	19.37	212	2972545	34.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3228269	34.74	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.84	94	36828	1.634 ng/ul#	88
45) Dimethylphthalate	13.73	163	168502	2.819 ng/ul	98

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

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