

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002004.D
 Acq On : 02 Mar 2020 23:48
 Operator : CG/JU
 Sample : L1616-13
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDQ7

Quant Time: Mar 03 03:56:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 03 03:51:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	286145	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1164088	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	734073	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1599687	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1574955	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1784346	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	31698	4.76	ng/uL	0.00
5) Phenol-d5	6.82	99	564121	25.84	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	321966	27.86	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	497161	26.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	455454	25.39	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	230959	25.83	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	251900	25.04	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	492417	25.41	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	652385	30.71	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1511076	27.20	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1797698	27.06	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	196473	18.50	ng/ul	0.00
58) Fluorene-d10	15.27	176	1336998	27.81	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	157992	17.92	ng/ul	0.00
71) Anthracene-d10	17.13	188	2009256	27.95	ng/ul	0.00
79) Pyrene-d10	19.37	212	2346328	28.46	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2562537	28.07	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.85	94	26048	1.150	ng/ul 97
45) Dimethylphthalate	13.73	163	125708	2.192	ng/ul 100

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

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