

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002012.D
 Acq On : 03 Mar 2020 05:26
 Operator : CG/JU
 Sample : L1616-21
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDR5

Quant Time: Mar 03 06:30:07 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	310237	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1267553	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	794264	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1688401	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1629075	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1865689	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.17	96	34121	4.73	ng/uL	0.00
5) Phenol-d5	6.82	99	629821	26.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	359560	28.69	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	548504	26.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	525406	27.02	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	258167	26.51	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	292626	26.71	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	548350	25.98	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	752935	32.55	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1657956	27.58	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1999150	27.81	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	228864	19.92	ng/ul	0.00
58) Fluorene-d10	15.27	176	1457968	28.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	174510	18.75	ng/ul	0.00
71) Anthracene-d10	17.13	188	2193987	28.91	ng/ul	0.00
79) Pyrene-d10	19.37	212	2512017	29.46	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2752020	28.83	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.85	94	29003	1.181	ng/ul 98
45) Dimethylphthalate	13.73	163	133946	2.159	ng/ul 98

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

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