

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002003.D
 Acq On : 02 Mar 2020 23:14
 Operator : CG/JU
 Sample : L1616-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDQ6

Quant Time: Mar 03 03:56:06 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	281550	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1160084	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	724828	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1580807	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1589160	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1784794	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.17	96	44289	6.76	ng/uL	0.00
5) Phenol-d5	6.82	99	796390	37.08	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	456651	40.15	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	702205	37.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	650920	36.88	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	340337	38.19	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	373259	37.23	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	697120	36.09	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	882200	41.67	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	2086170	38.03	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2518325	38.39	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	292544	27.90	ng/ul	0.00
58) Fluorene-d10	15.27	176	1826978	38.49	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	231719	26.59	ng/ul	0.00
71) Anthracene-d10	17.13	188	2745518	38.64	ng/ul	0.00
79) Pyrene-d10	19.37	212	3172087	38.13	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3526703	38.62	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.85	94	38306	1.718	ng/ul	96
45) Dimethylphthalate	13.73	163	152349	2.691	ng/ul	99

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

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