

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP001992.D
 Acq On : 02 Mar 2020 16:27
 Operator : CG/JU
 Sample : L1616-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDP8

Quant Time: Mar 02 16:58:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 02 15:51:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	258050	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1060786	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	670926	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1419657	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1409276	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1584692	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.17	96	31380	5.23	ng/uL	0.00
5) Phenol-d5	6.82	99	563755	28.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	340169	32.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	507625	29.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	442583	27.36	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	231905	28.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	252208	27.51	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	493357	27.94	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	558202	28.83	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1575930	31.04	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1873595	30.85	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	178632	18.40	ng/ul	0.00
58) Fluorene-d10	15.27	176	1409664	32.09	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	157459	20.12	ng/ul	0.00
71) Anthracene-d10	17.13	188	2122496	33.27	ng/ul	0.00
79) Pyrene-d10	19.37	212	2561641	34.72	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2756042	33.99	ng/ul	0.00
Target Compounds						
6) Phenol	6.85	94	21437	1.049	ng/ul	93
45) Dimethylphthalate	13.73	163	143225	2.733	ng/ul	100

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

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