

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002008.D
 Acq On : 03 Mar 2020 03:11
 Operator : CG/JU
 Sample : L1616-17
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDR1

Quant Time: Mar 03 05:33:13 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	282732	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1156402	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	733351	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1591664	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1602319	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1811869	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.16	96	31158	4.74	ng/uL	0.00
5) Phenol-d5	6.82	99	552938	25.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	325937	28.54	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	498261	26.71	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	391082	22.07	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	233189	26.25	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	258805	25.90	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	490159	25.46	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	628738	29.79	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1583819	28.54	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1845317	27.80	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	204563	19.28	ng/ul	0.00
58) Fluorene-d10	15.27	176	1387884	28.90	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	166627	18.99	ng/ul	0.00
71) Anthracene-d10	17.13	188	2160289	30.20	ng/ul	0.00
79) Pyrene-d10	19.37	212	2566120	30.59	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2822914	30.45	ng/ul	0.00
Target Compounds						
6) Phenol	6.85	94	27149	1.213	ng/ul	95
45) Dimethylphthalate	13.73	163	150095	2.620	ng/ul	99

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

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