

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP001987.D
 Acq On : 02 Mar 2020 13:34
 Operator : CG/JU
 Sample : L1616-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDP7

Quant Time: Mar 02 16:00:09 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	296341	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1202492	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	755830	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1647401	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1646441	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1886715	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.17	96	42455	6.16	ng/uL	0.00
5) Phenol-d5	6.82	99	700847	31.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	424232	35.44	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	634368	32.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	559666	30.13	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	294209	31.85	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	313575	30.17	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	604818	30.21	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	818776	37.31	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1914090	33.46	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2302524	33.66	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	217550	19.89	ng/ul	0.00
58) Fluorene-d10	15.27	176	1723830	34.83	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	189136	20.83	ng/ul	0.00
71) Anthracene-d10	17.13	188	2581849	34.87	ng/ul	0.00
79) Pyrene-d10	19.37	212	2987819	34.67	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3255985	33.73	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.85	94	41447	1.766	ng/ul	96
45) Dimethylphthalate	13.74	163	247683	4.195	ng/ul	98

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

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