

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
 Data File : BP002037.D
 Acq On : 04 Mar 2020 00:07
 Operator : CG/JU
 Sample : L1617-17
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Mar 04 05:16:40 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	284748	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1154275	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	738921	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1604311	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1600095	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1817221	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	133	0.02	ng/uL	0.01
5) Phenol-d5	6.83	99	1565	0.07	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.98	67	1741	0.15	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	1878	0.10	ng/ul	0.01
13) 4-Methylphenol-d8	8.36	113	692	0.04	ng/ul	0.02
19) Nitrobenzene-d5	8.78	128	201	0.02	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	490	0.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	912	0.05	ng/ul	0.04
29) 4-Chloroaniline-d4	10.53	131	84	0.00	ng/ul	-0.01
44) Dimethylphthalate-d6	13.70	166	7211	0.13	ng/ul	0.02
47) Acenaphthylene-d8	13.97	160	7118	0.11	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	11	0.00	ng/ul	0.00
58) Fluorene-d10	15.28	176	6414	0.13	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	8	0.00	ng/ul	0.00
71) Anthracene-d10	17.13	188	20730	0.29	ng/ul	0.00
79) Pyrene-d10	19.37	212	12336	0.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	12431	0.13	ng/ul	0.00

Target Compounds Ovalue

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

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