

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030520\
 Data File : BP002082.D
 Acq On : 05 Mar 2020 22:11
 Operator : CG/JU
 Sample : L1780-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0DG3

Quant Time: Mar 06 06:33:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 05 06:21:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	295448	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1211364	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	775278	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1676668	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1659044	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1850290	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	33585	4.89	ng/uL	0.00
5) Phenol-d5	6.82	99	120639	5.35	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	310843	26.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	399999	20.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	241552	13.04	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	229629	24.68	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	245196	23.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	475302	23.57	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	35625	1.61	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1765744	30.09	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1963051	27.98	ng/ul	0.00
52) 4-Nitrophenol-d4	14.48	143	40794	3.64	ng/ul	0.00
58) Fluorene-d10	15.27	176	1588129	31.28	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	194510	21.04	ng/ul	0.00
71) Anthracene-d10	17.12	188	2599692	34.50	ng/ul	0.00
79) Pyrene-d10	19.37	212	3006027	34.61	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	3293333	34.79	ng/ul	-0.01

Target Compounds

					Ovalue
33) 4-Chloro-3-methylphenol	11.70	107	23657	1.187 ng/ul	96
45) Dimethylphthalate	13.73	163	100183	1.654 ng/ul	100

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

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