

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030920\
 Data File : BP002131.D
 Acq On : 09 Mar 2020 21:12
 Operator : CG/JU
 Sample : L1834-14
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0DK7

Quant Time: Mar 09 23:28:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 09 21:43:02 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	281322	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	1152213	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	739274	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1638205	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1651113	20.00	ng/ul	0.00
86) Perylene-d12	23.32	264	1871276	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	33679	5.15	ng/uL	0.00
5) Phenol-d5	6.82	99	125644	5.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	383527	33.75	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	462020	24.89	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	197214	11.18	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	281976	31.86	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	312567	31.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	536811	27.98	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	28478	1.35	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1917894	34.28	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	2274327	33.99	ng/ul	0.00
52) 4-Nitrophenol-d4	14.48	143	37547	3.51	ng/ul	0.00
58) Fluorene-d10	15.27	176	1707914	35.28	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	230409	25.51	ng/ul	0.00
71) Anthracene-d10	17.12	188	2513415	34.14	ng/ul	0.00
79) Pyrene-d10	19.37	212	3015313	34.89	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.18	264	3475645	36.30	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.84	94	87094	3.909 ng/ul	98
45) Dimethylphthalate	13.73	163	164761	2.853 ng/ul	99

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

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