

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021721\
 Data File : BP004792.D
 Acq On : 18 Feb 2021 02:27
 Operator : CG/JU
 Sample : M1379-14
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBGZ4

Quant Time: Feb 18 03:06:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 17 23:17:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	55849	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	201155	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	117303	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	238690	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	244173	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	260724	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	7012	4.85	ng/uL	0.00
5) Phenol-d5	6.95	99	123386	27.53	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.10	67	64918	27.89	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	105681	31.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	96666	27.41	ng/ul	0.02
19) Nitrobenzene-d5	8.93	128	47554	34.24	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	56186	36.86	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	94728	30.88	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	122361	35.73	ng/ul	0.00
44) Dimethylphthalate-d6	13.83	166	278601	32.72	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	347741	33.57	ng/ul	0.00
52) 4-Nitrophenol-d4	14.64	143	39590	26.60	ng/ul	0.03
58) Fluorene-d10	15.41	176	231405	31.71	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	39229	27.19	ng/ul	0.00
71) Anthracene-d10	17.27	188	353142	33.55	ng/ul	0.00
79) Pyrene-d10	19.50	212	392537	34.34	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.41	264	444990	33.74	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.98	94	8608	1.842	ng/ul 96
45) Dimethylphthalate	13.87	163	24752	2.819	ng/ul# 96

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

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