

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\
 Data File : BM024589.D
 Acq On : 22 Jan 2020 19:33
 Operator : JU
 Sample : L1118-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GASK4

Quant Time: Jan 23 01:25:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 23 01:22:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	234463	20.00	ng/ul	0.00
18) Naphthalene-d8	10.21	136	1022273	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	726775	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1681673	20.00	ng/ul	0.00
78) Chrysene-d12	21.06	240	1554874	20.00	ng/ul	0.00
86) Perylene-d12	23.17	264	1422824	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	19408	4.05	ng/uL	0.00
5) Phenol-d5	6.64	99	382064	21.98	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.80	67	223636	23.17	ng/ul	0.00
9) 2-Chlorophenol-d4	6.99	132	347178	23.39	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	344337	21.49	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	181570	25.28	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	213231	27.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.84	165	391519	22.36	ng/ul	0.00
29) 4-Chloroaniline-d4	10.35	131	388159	19.94	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	1291011	22.67	ng/ul	0.00
47) Acenaphthylene-d8	13.79	160	1599773	24.41	ng/ul	0.00
52) 4-Nitrophenol-d4	14.31	143	194821	20.50	ng/ul	0.00
58) Fluorene-d10	15.10	176	1187162	23.68	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.22	200	175319	18.26	ng/ul	0.00
71) Anthracene-d10	16.95	188	1890592	24.18	ng/ul	0.00
79) Pyrene-d10	19.25	212	2191246	26.37	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.04	264	1765183	23.92	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.67	94	48384	2.740 ng/ul	93
45) Dimethylphthalate	13.57	163	206404	3.566 ng/ul	100

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

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