

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM010920\
 Data File : BM024428.D
 Acq On : 09 Jan 2020 17:19
 Operator : JU
 Sample : K6466-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0BQ6

Quant Time: Jan 09 18:04:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM010820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 09 01:07:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	320870	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1263644	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	749678	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	1621463	20.00	ng/ul	0.00
78) Chrysene-d12	21.09	240	1575234	20.00	ng/ul	-0.01
86) Perylene-d12	23.23	264	1822556	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.10	96	31558	4.24	ng/uL	0.00
5) Phenol-d5	6.68	99	724046	28.90	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.84	67	474102	32.26	ng/ul	0.00
9) 2-Chlorophenol-d4	7.03	132	645244	30.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	588612	29.25	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	326570	34.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	293063	32.26	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.88	165	600768	30.14	ng/ul	0.00
29) 4-Chloroaniline-d4	10.39	131	875114	38.93	ng/ul	0.00
44) Dimethylphthalate-d6	13.56	166	2027338	37.65	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	2524816	36.32	ng/ul	0.00
52) 4-Nitrophenol-d4	14.35	143	293646	30.92	ng/ul	0.00
58) Fluorene-d10	15.14	176	1675168	35.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	181104	24.93	ng/ul	0.00
71) Anthracene-d10	16.99	188	2679092	35.28	ng/ul	0.00
79) Pyrene-d10	19.29	212	2965217	35.23	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.10	264	3109037	33.41	ng/ul	-0.01

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

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