

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022621\
 Data File : BM028901.D
 Acq On : 25 Feb 2021 17:49
 Operator : CG/JU
 Sample : M1482-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBGMS

Quant Time: Feb 25 23:24:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 25 15:24:23 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	15135	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	59347	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	32413	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	59499	20.00	ng/ul	0.00
78) Chrysene-d12	21.36	240	52985	20.00	ng/ul	0.00
86) Perylene-d12	23.57	264	54203	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	1538	4.84	ng/uL	0.00
5) Phenol-d5	6.99	99	7149	6.20	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	20270	31.20	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	26173	26.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	14126	14.60	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	14462	34.43	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	16399	34.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	25953	28.62	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	1556	1.45	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	77011	32.96	ng/ul	0.00
47) Acenaphthylene-d8	14.16	160	102430	34.42	ng/ul	0.00
52) 4-Nitrophenol-d4	14.71	143	1722	4.26	ng/ul	0.00
58) Fluorene-d10	15.46	176	68231	34.73	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.60	200	11001	30.23	ng/ul	0.00
71) Anthracene-d10	17.30	188	108997	40.22	ng/ul	0.00
79) Pyrene-d10	19.59	212	109266	41.37	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.44	264	114184	40.87	ng/ul	0.00

Target Compounds Ovalue

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

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