

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040921\
 Data File : BM029442.D
 Acq On : 09 Apr 2021 21:40
 Operator : CG/JU
 Sample : M1881-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ESQF8

Quant Time: Apr 10 01:18:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 09 11:03:41 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	110023	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	433281	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	229939	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	399443	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	338491	20.00	ng/ul	0.00
86) Perylene-d12	23.45	264	365799	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	5681	1.97	ng/uL	0.00
5) Phenol-d5	6.86	99	89213	9.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	66772	11.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.22	132	75137	10.78	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	31115	4.38	ng/ul	0.00
19) Nitrobenzene-d5	8.84	128	34582	11.53	ng/ul	0.00
22) 2-Nitrophenol-d4	9.56	143	35046	11.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	58176	9.06	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	66733	8.54	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	191257	11.61	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	233030	11.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	19198	6.36	ng/ul	0.00
58) Fluorene-d10	15.32	176	152097	10.93	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	10890	4.80	ng/ul	0.00
71) Anthracene-d10	17.16	188	214461	11.81	ng/ul	0.00
79) Pyrene-d10	19.45	212	238374	14.67	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.31	264	248499	13.31	ng/ul	0.00

Target Compounds

					Ovalue
4) Benzaldehyde	6.84	77	7012	1.374	ng/ul 92
6) Phenol	6.89	94	19271	1.960	ng/ul 93
14) Acetophenone	8.51	105	39613	3.411	ng/ul 95

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

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