

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
 Data File : BM030860.D
 Acq On : 07 Jul 2021 17:07
 Operator : CG/JU
 Sample : M2854-14
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGD8

Quant Time: Jul 07 17:57:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 06 14:32:50 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	84558	20.000	ng/ul	0.00
20) Naphthalene-d8	10.545	136	333186	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.392	164	195787	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.133	188	364315	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	357543	20.000	ng/ul	0.00
88) Perylene-d12	23.551	264	396139	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	5440	2.514	ng/uL	0.00
4) Pyridine-d5	3.699	84	23791	4.204	ng/ul	0.02
7) Phenol-d5	6.934	99	81987	11.096	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	60180	12.645	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	69702	12.296	ng/ul	0.00
15) 4-Methylphenol-d8	8.463	113	64190	10.841	ng/ul	0.00
21) Nitrobenzene-d5	8.922	128	31814	12.838	ng/ul	0.00
24) 2-Nitrophenol-d4	9.640	143	31720	11.699	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	58323	11.177	ng/ul	0.00
31) 4-Chloroaniline-d4	10.693	131	74717	10.083	ng/ul	0.00
46) Dimethylphthalate-d6	13.804	166	182682	13.328	ng/ul	0.00
49) Acenaphthylene-d8	14.086	160	213813	12.349	ng/ul	0.00
54) 4-Nitrophenol-d4	14.616	143	15211	6.377	ng/ul	0.02
60) Fluorene-d10	15.386	176	155751	12.904	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	14263	6.106	ng/ul	0.00
73) Anthracene-d10	17.227	188	234305	13.801	ng/ul	0.00
81) Pyrene-d10	19.521	212	240409	12.192	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.409	264	210662	10.302	ng/ul	0.00

Target Compounds Qvalue

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

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