

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM063021\
 Data File : BM030745.D
 Acq On : 01 Jul 2021 16:14
 Operator : CG/JU
 Sample : M2825-17
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDR5

Manual Integrations
 APPROVED

mohammad
 7/2/2021 12:02:47 PM

Quant Time: Jul 01 16:46:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 01 05:56:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	75643	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	278648	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.404	164	155914	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.145	188	300852	20.000	ng/u1	0.00
79) Chrysene-d12	21.309	240	320723	20.000	ng/u1	0.00
88) Perylene-d12	23.562	264	375658	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	1434	0.769	ng/uL	0.02
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	6.946	99	16156	2.476	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	11993	2.925	ng/u1	0.00
11) 2-Chlorophenol-d4	7.310	132	13668	2.714	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	11204	2.207	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	5351	2.573	ng/u1	0.00
24) 2-Nitrophenol-d4	9.657	143	4681	2.026	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	9031	2.040	ng/u1	0.02
31) 4-Chloroaniline-d4	10.728	131	11681m	1.892	ng/u1	0.02
46) Dimethylphthalate-d6	13.822	166	29536	2.744	ng/u1	0.00
49) Acenaphthylene-d8	14.098	160	37083	2.654	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	861m	0.422	ng/u1	0.06
60) Fluorene-d10	15.398	176	28194	2.953	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	1041m	0.529	ng/u1	0.01
73) Anthracene-d10	17.245	188	43101	3.070	ng/u1	0.00
81) Pyrene-d10	19.527	212	46211	2.750	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.415	264	52146	2.670	ng/u1	-0.01

Target Compounds Qvalue

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

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