

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
 Data File : BM032177.D
 Acq On : 26 Sep 2021 05:41
 Operator : CG/JU
 Sample : M3739-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B73

Quant Time: Sep 26 06:48:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM092421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 25 01:15:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	380648	20.000	ng/u1	0.00
20) Naphthalene-d8	10.454	136	1463181	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.307	164	824416	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.036	188	1566788	20.000	ng/u1	0.00
79) Chrysene-d12	21.194	240	1421082	20.000	ng/u1	0.00
88) Perylene-d12	23.359	264	1433202	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.990	96	47451	5.264	ng/uL	0.00
4) Pyridine-d5	3.419	84	210655	8.231	ng/u1	0.00
7) Phenol-d5	6.842	99	192576	6.374	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.984	67	499042	27.468	ng/u1	0.00
11) 2-Chlorophenol-d4	7.201	132	539016	22.978	ng/u1	0.00
15) 4-Methylphenol-d8	8.384	113	344001	14.560	ng/u1	-0.01
21) Nitrobenzene-d5	8.819	128	290656	33.259	ng/u1	0.00
24) 2-Nitrophenol-d4	9.542	143	291792	31.576	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.083	165	556639	26.350	ng/u1	0.00
31) 4-Chloroaniline-d4	10.589	131	777910	27.246	ng/u1	0.00
46) Dimethylphthalate-d6	13.713	166	1769871	33.042	ng/u1	0.00
49) Acenaphthylene-d8	14.001	160	2102575	30.896	ng/u1	0.00
54) 4-Nitrophenol-d4	14.501	143	45862	5.263	ng/u1	-0.01
60) Fluorene-d10	15.295	176	1567079	34.286	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.407	200	196904	26.592	ng/u1	-0.01
73) Anthracene-d10	17.130	188	2546394	38.951	ng/u1	0.00
81) Pyrene-d10	19.418	212	2763677	36.962	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.224	264	2552793	37.954	ng/u1	-0.01

Target Compounds Qvalue

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

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