

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037332.D
 Acq On : 03 Nov 2022 00:38
 Operator : CG/JU
 Sample : N5321-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWW1

Quant Time: Nov 03 01:21:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 02 03:33:37 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.851	152	318201	20.000 ng/u1	0.00
20) Naphthalene-d8	10.651	136	1399319	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	920885	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1805945	20.000 ng/u1	0.00
79) Chrysene-d12	21.397	240	1398038	20.000 ng/u1	0.00
88) Perylene-d12	23.744	264	1282124	20.000 ng/u1	-0.01

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.275	96	37873	5.247 ng/uL	0.00
4) Pyridine-d5	3.699	84	243137	11.284 ng/u1	0.00
7) Phenol-d5	7.028	99	200643	7.245 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	596694	35.114 ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	588786	27.828 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	296817	13.126 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	394217	37.714 ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	425490	37.370 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	701541	33.420 ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	471536	14.699 ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2601538	40.859 ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	3034436	39.641 ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	69342	5.448 ng/u1	0.00
60) Fluorene-d10	15.480	176	2354042	41.459 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	399533	34.929 ng/u1	0.00
73) Anthracene-d10	17.327	188	3260315	39.200 ng/u1	0.00
81) Pyrene-d10	19.615	212	3476702	47.590 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.591	264	2788124	42.246 ng/u1	-0.01

Target Compounds	Qvalue

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

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