

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037333.D
 Acq On : 03 Nov 2022 01:14
 Operator : CG/JU
 Sample : N5321-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWW6

Quant Time: Nov 03 02:51:39 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	381071	20.000 ng/u1	0.00
20) Naphthalene-d8	10.651	136	1643672	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	1040669	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	2068647	20.000 ng/u1	0.00
79) Chrysene-d12	21.398	240	1722746	20.000 ng/u1	0.00
88) Perylene-d12	23.744	264	1496006	20.000 ng/u1	-0.01

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.275	96	29383	3.399 ng/uL	0.00
4) Pyridine-d5	3.699	84	252428	9.782 ng/u1	0.00
7) Phenol-d5	7.028	99	170985	5.156 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	521469	25.624 ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	517051	20.406 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	319157	11.785 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	340783	27.756 ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	355351	26.570 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	637493	25.854 ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	671488	17.820 ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2346185	32.607 ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2680728	30.989 ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	63262	4.399 ng/u1	0.00
60) Fluorene-d10	15.480	176	2115739	32.973 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	356122	27.180 ng/u1	0.00
73) Anthracene-d10	17.327	188	3326151	34.913 ng/u1	0.00
81) Pyrene-d10	19.615	212	3588315	39.860 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.592	264	2752206	35.739 ng/u1	-0.01

Target Compounds	Qvalue

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

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