

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037326.D
 Acq On : 02 Nov 2022 21:05
 Operator : CG/JU
 Sample : N5301-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWM4

Quant Time: Nov 02 22:55:36 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	7.857	152	331493	20.000 ng/u1	0.00
20) Naphthalene-d8	10.657	136	1503664	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	996526	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	2087835	20.000 ng/u1	0.00
79) Chrysene-d12	21.397	240	1992070	20.000 ng/u1	0.00
88) Perylene-d12	23.744	264	1732721	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.275	96	31700	4.215 ng/uL	0.00
4) Pyridine-d5	3.699	84	214409	9.551 ng/u1	0.00
7) Phenol-d5	7.028	99	200383	6.946 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	582898	32.927 ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	579851	26.307 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	390972	16.596 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	379218	33.762 ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	398381	32.561 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	689834	30.582 ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	868708	25.201 ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2433478	35.318 ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2885896	34.839 ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	67376	4.892 ng/u1	0.00
60) Fluorene-d10	15.480	176	2247918	36.585 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	366814	27.739 ng/u1	0.00
73) Anthracene-d10	17.327	188	3582448	37.257 ng/u1	0.00
81) Pyrene-d10	19.615	212	4075383	39.150 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.597	264	3356696	37.634 ng/u1	0.00

Target Compounds Qvalue

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

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