

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037354.D
 Acq On : 03 Nov 2022 19:40
 Operator : CG/JU
 Sample : N5276-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4W8

Quant Time: Nov 04 01:05:49 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	338842	20.000 ng/u1	0.00
20) Naphthalene-d8	10.651	136	1488023	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	954234	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1855363	20.000 ng/u1	0.00
79) Chrysene-d12	21.397	240	1385504	20.000 ng/u1	0.00
88) Perylene-d12	23.744	264	1231367	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.269	96	31268	4.068 ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000 ng/u1	
7) Phenol-d5	7.028	99	215738	7.316 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	632843	34.972 ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	633439	28.115 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	402560	16.718 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	409117	36.806 ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	430324	35.541 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	699653	31.343 ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	22329	0.655 ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2423879	36.738 ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2921951	36.837 ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	63505	4.815 ng/u1	0.00
60) Fluorene-d10	15.474	176	2257718	38.373 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	340341	28.962 ng/u1	0.00
73) Anthracene-d10	17.327	188	3412926	39.941 ng/u1	0.00
81) Pyrene-d10	19.615	212	3491667	48.227 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.591	264	2550370	40.236 ng/u1	-0.01

Target Compounds	Qvalue

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

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