

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037353.D
 Acq On : 03 Nov 2022 19:04
 Operator : CG/JU
 Sample : N5276-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4W3

Quant Time: Nov 04 01:05:31 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	333200	20.000 ng/u1	0.00
20) Naphthalene-d8	10.651	136	1512647	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	1009105	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	2120669	20.000 ng/u1	0.00
79) Chrysene-d12	21.397	240	2035833	20.000 ng/u1	0.00
88) Perylene-d12	23.744	264	1811525	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.269	96	26761	3.540 ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000 ng/u1	
7) Phenol-d5	7.028	99	184526	6.363 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	503235	28.281 ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	500589	22.594 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	338947	14.314 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	331238	29.315 ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	344402	27.982 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	581867	25.642 ng/u1	0.00
31) 4-Chloroaniline-d4	10.816	131	37929	1.094 ng/u1	0.01
46) Dimethylphthalate-d6	13.904	166	2188168	31.362 ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2549485	30.394 ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	65508	4.697 ng/u1	0.00
60) Fluorene-d10	15.480	176	2071745	33.297 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	343550	25.577 ng/u1	0.00
73) Anthracene-d10	17.327	188	3536359	36.208 ng/u1	0.00
81) Pyrene-d10	19.615	212	3975546	37.370 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.591	264	3410316	36.572 ng/u1	-0.01

Target Compounds Qvalue

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

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