

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037341.D
 Acq On : 03 Nov 2022 11:48
 Operator : CG/JU
 Sample : N5276-20
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4Y0

Quant Time: Nov 04 01:00:01 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	256874	20.000 ng/u1	0.00
20) Naphthalene-d8	10.651	136	1148984	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	738823	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1531448	20.000 ng/u1	0.00
79) Chrysene-d12	21.397	240	1447165	20.000 ng/u1	0.00
88) Perylene-d12	23.744	264	1311296	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.269	96	32370	5.555 ng/uL	0.00
4) Pyridine-d5	3.704	84	94221	5.417 ng/u1	0.00
7) Phenol-d5	7.022	99	204640	9.154 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	599697	43.716 ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	586707	34.350 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	390823	21.409 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	381618	44.463 ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	389461	41.658 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	659446	38.259 ng/u1	0.00
31) 4-Chloroaniline-d4	10.798	131	777774	29.528 ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2369789	46.390 ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2854813	46.484 ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	67575	6.618 ng/u1	0.00
60) Fluorene-d10	15.474	176	2277332	49.991 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	351263	36.213 ng/u1	0.00
73) Anthracene-d10	17.321	188	3608123	51.157 ng/u1	-0.01
81) Pyrene-d10	19.615	212	4070088	53.821 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.591	264	3478392	51.532 ng/u1	-0.01

Target Compounds	Qvalue

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

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