

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037357.D
 Acq On : 03 Nov 2022 21:27
 Operator : CG/JU
 Sample : N5276-11
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4X1

Quant Time: Nov 04 01:06:46 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	320426	20.000	ng/u1	0.00
20) Naphthalene-d8	10.651	136	1450158	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.486	164	947932	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1954184	20.000	ng/u1	0.00
79) Chrysene-d12	21.397	240	1848503	20.000	ng/u1	0.00
88) Perylene-d12	23.744	264	1615318	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	30547	4.202	ng/uL	0.00
4) Pyridine-d5	3.693	84	174	0.008	ng/u1	0.00
7) Phenol-d5	7.022	99	185566	6.654	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.186	67	536751	31.367	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	532082	24.973	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	353174	15.510	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	348280	32.151	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.739	143	356810	30.239	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	596410	27.415	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	45899	1.381	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2219811	33.869	ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2577408	32.710	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	64310	4.909	ng/u1	0.00
60) Fluorene-d10	15.474	176	2110745	36.113	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	350093	28.285	ng/u1	0.00
73) Anthracene-d10	17.321	188	3525627	39.174	ng/u1	-0.01
81) Pyrene-d10	19.615	212	3984267	41.247	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.591	264	3339146	40.158	ng/u1	-0.01

Target Compounds Qvalue

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

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