

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037366.D
 Acq On : 04 Nov 2022 03:20
 Operator : CG/JU
 Sample : N5276-17
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4X7

Quant Time: Nov 04 03:59:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:39:09 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	278666	20.000	ng/u1	0.00
20) Naphthalene-d8	10.645	136	1205456	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.480	164	726965	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	1353808	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1044515	20.000	ng/u1	0.00
88) Perylene-d12	23.733	264	1033431	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	37564	5.942	ng/uL	0.00
4) Pyridine-d5	3.699	84	180232	9.551	ng/u1	0.00
7) Phenol-d5	7.022	99	251104	10.354	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	683412	45.923	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	680562	36.729	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	449260	22.686	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	435796	48.397	ng/u1	0.00
24) 2-Nitrophenol-d4	9.734	143	438777	44.734	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	751155	41.538	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	891895	32.274	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2514116	50.018	ng/u1	0.00
49) Acenaphthylene-d8	14.175	160	3066566	50.747	ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	69221	6.890	ng/u1	0.00
60) Fluorene-d10	15.474	176	2339996	52.205	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	333836	38.932	ng/u1	0.00
73) Anthracene-d10	17.321	188	3447801	55.298	ng/u1	0.00
81) Pyrene-d10	19.609	212	3429795	62.837	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.586	264	2993330	56.269	ng/u1	0.00

Target Compounds Qvalue

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

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