

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037365.D
 Acq On : 04 Nov 2022 02:45
 Operator : CG/JU
 Sample : N5276-09
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4W7

Quant Time: Nov 04 03:39:48 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	337072	20.000	ng/u1	0.00
20) Naphthalene-d8	10.645	136	1492653	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.480	164	928823	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	1836354	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1502874	20.000	ng/u1	0.00
88) Perylene-d12	23.733	264	1313613	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	32404	4.238	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.022	99	213486	7.278	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	614950	34.162	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	607236	27.093	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	396883	16.568	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	394329	35.366	ng/u1	0.00
24) 2-Nitrophenol-d4	9.734	143	400356	32.964	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	681536	30.437	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	38626	1.129	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	2290913	35.672	ng/u1	0.00
49) Acenaphthylene-d8	14.174	160	2782706	36.042	ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	61310	4.776	ng/u1	0.00
60) Fluorene-d10	15.474	176	2158008	37.682	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	324625	27.910	ng/u1	0.00
73) Anthracene-d10	17.321	188	3362136	39.754	ng/u1	0.00
81) Pyrene-d10	19.609	212	3598361	45.819	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.586	264	2741348	40.541	ng/u1	0.00

Target Compounds Qvalue

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

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