

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
 Data File : BM044367.D
 Acq On : 27 Feb 2024 15:16
 Operator : MA/JU
 Sample : P1469-10
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Feb 27 22:28:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 27 00:07:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	177121	20.000	ng/ul	0.01
20) Naphthalene-d8	10.746	136	755724	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.581	164	434713	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.327	188	900475	20.000	ng/ul	0.00
79) Chrysene-d12	21.527	240	776510	20.000	ng/ul	0.00
88) Perylene-d12	23.945	264	854418	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	27491	4.897	ng/uL	0.00
4) Pyridine-d5	3.781	84	375960	24.522	ng/ul	0.00
7) Phenol-d5	7.105	99	457180	26.944	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	292970	26.664	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	362743	28.385	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	356914	27.800	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	178220	29.004	ng/ul	0.01
24) 2-Nitrophenol-d4	9.834	143	195084	31.119	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.363	165	324900	29.380	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	517694	27.874	ng/ul	0.01
46) Dimethylphthalate-d6	14.004	166	1072930	32.002	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1246707	32.320	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	159686	23.936	ng/ul	0.02
60) Fluorene-d10	15.575	176	941475	33.177	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	117577	21.044	ng/ul	0.00
73) Anthracene-d10	17.427	188	1451984	33.817	ng/ul	0.00
81) Pyrene-d10	19.727	212	1653271	35.390	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	1527631	34.892	ng/ul	0.02

Target Compounds	Qvalue

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

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