

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
 Data File : BM044366.D
 Acq On : 27 Feb 2024 14:40
 Operator : MA/JU
 Sample : P1469-06
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Feb 27 22:27:54 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	194533	20.000	ng/ul	0.01
20) Naphthalene-d8	10.745	136	816605	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.580	164	465228	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	963476	20.000	ng/ul	0.01
79) Chrysene-d12	21.527	240	821705	20.000	ng/ul	0.00
88) Perylene-d12	23.944	264	900844	20.000	ng/ul	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	32800	5.320	ng/uL	0.00
4) Pyridine-d5	3.787	84	147897	8.783	ng/ul	0.01
7) Phenol-d5	7.104	99	567907	30.474	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	356615	29.551	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	443617	31.607	ng/ul	0.00
15) 4-Methylphenol-d8	8.651	113	424071	30.075	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	220263	33.174	ng/ul	0.01
24) 2-Nitrophenol-d4	9.828	143	222801	32.890	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	379742	31.779	ng/ul	0.00
31) 4-Chloroaniline-d4	10.892	131	577554	28.778	ng/ul	0.01
46) Dimethylphthalate-d6	14.004	166	1184951	33.025	ng/ul	0.00
49) Acenaphthylene-d8	14.274	160	1441467	34.918	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	163394	22.885	ng/ul	0.02
60) Fluorene-d10	15.574	176	1069895	35.230	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	108514	18.152	ng/ul	0.00
73) Anthracene-d10	17.427	188	1645758	35.824	ng/ul	0.00
81) Pyrene-d10	19.727	212	1833791	37.095	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.791	264	1663474	36.037	ng/ul	0.02

Target Compounds	Qvalue

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

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