

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048271.D
 Acq On : 29 Oct 2024 01:48
 Operator : RC/JU
 Sample : P4529-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F7H23

Quant Time: Oct 29 02:34:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Oct 27 23:39:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	229954	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	961544	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	601575	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1139558	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	956631	20.000	ng/u1	0.00
88) Perylene-d12	24.268	264	1061048	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	29265	5.010	ng/uL	0.00
4) Pyridine-d5	3.587	84	396418	23.415	ng/u1	0.00
7) Phenol-d5	6.893	99	512397	25.937	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	341830	27.932	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	434830	27.066	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	437809	27.603	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	214242	27.177	ng/u1	0.00
24) 2-Nitrophenol-d4	9.598	143	240421	28.152	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	409803	26.296	ng/u1	0.00
31) 4-Chloroaniline-d4	10.645	131	634383	29.316	ng/u1	0.00
46) Dimethylphthalate-d6	13.774	166	1397558	31.087	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	1571194	30.996	ng/u1	0.00
54) 4-Nitrophenol-d4	14.586	143	142127	17.288	ng/u1	0.00
60) Fluorene-d10	15.357	176	1231804	32.427	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	165678	22.000	ng/u1	0.00
73) Anthracene-d10	17.204	188	1863091	34.575	ng/u1	0.00
81) Pyrene-d10	19.492	212	2031422	34.994	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.068	264	1952330	34.885	ng/u1	-0.01

Target Compounds	Qvalue

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

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