

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048292.D
 Acq On : 29 Oct 2024 18:04
 Operator : RC/JU
 Sample : PB164416BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK416

Quant Time: Oct 29 22:59:05 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	176168	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	734496	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	448269	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	844644	20.000	ng/u1	-0.01
79) Chrysene-d12	21.327	240	672013	20.000	ng/u1	0.00
88) Perylene-d12	24.268	264	748925	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	27540	6.155	ng/uL	0.00
4) Pyridine-d5	3.587	84	386407	29.792	ng/u1	0.00
7) Phenol-d5	6.892	99	442197	29.217	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.045	67	292200	31.167	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.251	132	376395	30.581	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	356550	29.343	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	181152	30.083	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	197156	30.222	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.133	165	330160	27.734	ng/u1	0.00
31) 4-Chloroaniline-d4	10.645	131	482232	29.173	ng/u1	0.00
46) Dimethylphthalate-d6	13.769	166	999834	29.846	ng/u1	-0.01
49) Acenaphthylene-d8	14.051	160	1155518	30.592	ng/u1	0.00
54) 4-Nitrophenol-d4	14.580	143	115313	18.823	ng/u1	0.00
60) Fluorene-d10	15.351	176	842449	29.762	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	118918	21.304	ng/u1	0.00
73) Anthracene-d10	17.198	188	1233572	30.886	ng/u1	-0.01
81) Pyrene-d10	19.492	212	1329786	32.609	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.068	264	1254143	31.749	ng/u1	-0.01

Target Compounds Qvalue

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

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