

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048293.D
 Acq On : 29 Oct 2024 18:43
 Operator : RC/JU
 Sample : PB164419BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK419

Quant Time: Oct 29 23:01:09 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	174382	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	715130	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	437331	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	836029	20.000	ng/u1	-0.01
79) Chrysene-d12	21.327	240	712546	20.000	ng/u1	0.00
88) Perylene-d12	24.268	264	793462	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	27201	6.141	ng/uL	0.00
4) Pyridine-d5	3.587	84	371717	28.953	ng/u1	0.00
7) Phenol-d5	6.892	99	420208	28.049	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	275110	29.645	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	359941	29.544	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	332797	27.669	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	170137	29.018	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	183670	28.917	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.133	165	308034	26.576	ng/u1	0.00
31) 4-Chloroaniline-d4	10.645	131	449325	27.918	ng/u1	0.00
46) Dimethylphthalate-d6	13.768	166	950372	29.079	ng/u1	-0.01
49) Acenaphthylene-d8	14.051	160	1083650	29.407	ng/u1	0.00
54) 4-Nitrophenol-d4	14.586	143	110782	18.536	ng/u1	0.00
60) Fluorene-d10	15.351	176	813645	29.463	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	115490	20.903	ng/u1	0.00
73) Anthracene-d10	17.198	188	1218763	30.829	ng/u1	-0.01
81) Pyrene-d10	19.492	212	1328832	30.732	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.068	264	1298367	31.024	ng/u1	-0.01

Target Compounds Qvalue

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

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