

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048264.D
 Acq On : 28 Oct 2024 21:13
 Operator : RC/JU
 Sample : PB164423BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK423

Quant Time: Oct 28 23:19:43 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	241779	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1014916	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	633615	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1224062	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1009226	20.000	ng/u1	0.00
88) Perylene-d12	24.279	264	1107309	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	32003	5.211	ng/uL	0.00
4) Pyridine-d5	3.587	84	450720	25.320	ng/u1	0.00
7) Phenol-d5	6.898	99	503571	24.243	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	332184	25.817	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	431672	25.555	ng/u1	0.00
15) 4-Methylphenol-d8	8.433	113	405427	24.311	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	212221	25.505	ng/u1	0.00
24) 2-Nitrophenol-d4	9.598	143	230209	25.539	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.139	165	371012	22.555	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	564419	24.711	ng/u1	0.00
46) Dimethylphthalate-d6	13.774	166	1235714	26.097	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	1377851	25.808	ng/u1	0.00
54) 4-Nitrophenol-d4	14.580	143	157763	18.219	ng/u1	0.00
60) Fluorene-d10	15.357	176	1046684	26.160	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	170169	21.036	ng/u1	0.00
73) Anthracene-d10	17.203	188	1578873	27.278	ng/u1	0.00
81) Pyrene-d10	19.492	212	1725117	28.168	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.074	264	1610630	27.577	ng/u1	0.00

Target Compounds Qvalue

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

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