

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048309.D
 Acq On : 30 Oct 2024 05:50
 Operator : RC/JU
 Sample : PB164514BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK514

Quant Time: Oct 30 06:28:52 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	209292	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	848161	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	510161	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	969108	20.000	ng/u1	-0.01
79) Chrysene-d12	21.321	240	811425	20.000	ng/u1	-0.01
88) Perylene-d12	24.262	264	899797	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	32178	6.053	ng/uL	0.00
4) Pyridine-d5	3.581	84	451001	29.269	ng/u1	0.00
7) Phenol-d5	6.893	99	515409	28.665	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.045	67	336246	30.189	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.251	132	439441	30.053	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	402975	27.915	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	209683	30.154	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	224556	29.809	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.133	165	389082	28.304	ng/u1	0.00
31) 4-Chloroaniline-d4	10.645	131	537000	28.133	ng/u1	0.00
46) Dimethylphthalate-d6	13.769	166	1114707	29.238	ng/u1	-0.01
49) Acenaphthylene-d8	14.051	160	1285422	29.903	ng/u1	0.00
54) 4-Nitrophenol-d4	14.580	143	128495	18.430	ng/u1	0.00
60) Fluorene-d10	15.351	176	941053	29.212	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	131096	20.469	ng/u1	0.00
73) Anthracene-d10	17.198	188	1422907	31.051	ng/u1	-0.01
81) Pyrene-d10	19.492	212	1565371	31.791	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.062	264	1530497	32.248	ng/u1	-0.02

Target Compounds Qvalue

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

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