

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070225\
 Data File : BM050338.D
 Acq On : 02 Jul 2025 10:41
 Operator : RC/JU
 Sample : PB168464BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK464

Quant Time: Jul 02 12:26:03 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM070125.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 01 15:14:38 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	376734	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	1393699	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	853516	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1599874	20.000	ng/u1	0.00
79) Chrysene-d12	21.445	240	1468111	20.000	ng/u1	0.00
88) Perylene-d12	24.480	264	1564041	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	58297	6.387	ng/uL	0.00
4) Pyridine-d5	3.710	84	883829	35.767	ng/u1	0.00
7) Phenol-d5	7.022	99	949847	34.777	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	605047	34.934	ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	768714	34.557	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	707859	34.019	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	305635	31.714	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	285100	31.738	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	690338	31.620	ng/u1	0.00
31) 4-Chloroaniline-d4	10.786	131	989706	34.121	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	1788726	30.769	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2232987	31.370	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	239596	22.932	ng/u1	0.00
60) Fluorene-d10	15.480	176	1692171	32.065	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	123093	17.813	ng/u1	0.00
73) Anthracene-d10	17.321	188	2561162	32.282	ng/u1	0.00
81) Pyrene-d10	19.603	212	2780370	35.452	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.280	264	2599911	32.788	ng/u1	0.00

Target Compounds	Qvalue

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

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