

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070225\
 Data File : BM050343.D
 Acq On : 02 Jul 2025 14:03
 Operator : RC/JU
 Sample : PB168471BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK471

Quant Time: Jul 02 15:01:11 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	335715	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	1256022	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	767611	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1482528	20.000	ng/u1	0.00
79) Chrysene-d12	21.450	240	1409899	20.000	ng/u1	0.00
88) Perylene-d12	24.485	264	1531736	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	52810	6.493	ng/uL	0.00
4) Pyridine-d5	3.710	84	842926	38.280	ng/u1	0.00
7) Phenol-d5	7.022	99	875282	35.963	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	550612	35.676	ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	694362	35.028	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	651298	35.126	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	270387	31.132	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	240608	29.721	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	609970	31.001	ng/u1	0.00
31) 4-Chloroaniline-d4	10.786	131	949827	36.336	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	1623067	31.044	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2021707	31.581	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	233532	24.853	ng/u1	0.00
60) Fluorene-d10	15.480	176	1541924	32.487	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	111935	17.481	ng/u1	0.00
73) Anthracene-d10	17.321	188	2412634	32.817	ng/u1	0.00
81) Pyrene-d10	19.603	212	2687850	35.687	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.280	264	2522960	32.489	ng/u1	0.00

Target Compounds	Qvalue

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

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