

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
 Data File : BM043307.D
 Acq On : 13 Dec 2023 18:37
 Operator : MA/JU
 Sample : PB157727BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK727

Quant Time: Dec 13 22:39:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 13 22:33:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	380492	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1840579	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.621	164	1093876	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.362	188	2103102	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.533	240	1491641	20.000	ng/u1	# 0.00
88) Perylene-d12	23.950	264	1403158	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	57616	5.943	ng/uL	0.00
4) Pyridine-d5	3.816	84	904416	31.441	ng/u1	0.00
7) Phenol-d5	7.157	99	1183455	31.329	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	738507	31.581	ng/u1	0.00
11) 2-Chlorophenol-d4	7.528	132	877946	30.487	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	944828	31.616	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	448928	28.616	ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	447569	29.586	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	835549	24.780	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	1339498	27.609	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2677435	27.922	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	3051999	27.511	ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	494533	28.782	ng/u1	0.00
60) Fluorene-d10	15.616	176	2219282	26.551	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	88422	7.639	ng/u1	0.00
73) Anthracene-d10	17.462	188	3495751	30.724	ng/u1	0.00
81) Pyrene-d10	19.745	212	3510554	36.782	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.797	264	2606987	31.841	ng/u1	0.00

Target Compounds Qvalue

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

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