

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110824\
 Data File : BM048536.D
 Acq On : 08 Nov 2024 18:54
 Operator : RC/JU
 Sample : PB164769BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK769

Quant Time: Nov 08 22:05:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 07 22:50:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	109143	20.000	ng/u1	0.00
20) Naphthalene-d8	10.451	136	449342	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	263640	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.056	188	541500	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	534286	20.000	ng/u1	0.00
88) Perylene-d12	24.197	264	614816	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	18623	6.802	ng/uL	0.00
4) Pyridine-d5	3.546	84	246076	32.614	ng/u1	0.00
7) Phenol-d5	6.845	99	292375	35.190	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	173946	34.856	ng/u1	0.00
11) 2-Chlorophenol-d4	7.198	132	250240	36.360	ng/u1	0.00
15) 4-Methylphenol-d8	8.380	113	228561	34.986	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	114502	35.158	ng/u1	0.00
24) 2-Nitrophenol-d4	9.545	143	131927	38.084	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.080	165	219160	34.056	ng/u1	0.00
31) 4-Chloroaniline-d4	10.598	131	306080	33.486	ng/u1	0.00
46) Dimethylphthalate-d6	13.727	166	630987	37.493	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	744232	37.508	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	85587	27.381	ng/u1	0.01
60) Fluorene-d10	15.309	176	550246	37.456	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	72977	26.790	ng/u1	0.00
73) Anthracene-d10	17.156	188	865400	37.766	ng/u1	0.00
81) Pyrene-d10	19.450	212	1052743	39.262	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.997	264	1131067	39.374	ng/u1	0.00

Target Compounds	Qvalue

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

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