

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062846.D
 Acq On : 15 Sep 2024 14:57
 Operator : JU/RC
 Sample : PB163064BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK064

Quant Time: Sep 15 23:35:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 01:28:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.897	152	33803	20.000	ng/u1	0.00
20) Naphthalene-d8	10.694	136	137098	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.531	164	116930	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.275	188	325883	20.000	ng/u1	0.00
79) Chrysene-d12	21.517	240	366681	20.000	ng/u1	0.00
88) Perylene-d12	24.584	264	408368	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.362	96	4764	6.371	ng/uL	0.00
4) Pyridine-d5	3.767	84	68904	28.947	ng/u1	0.00
7) Phenol-d5	7.069	99	84529	28.517	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	50641	28.496	ng/u1	0.00
11) 2-Chlorophenol-d4	7.433	132	65669	30.415	ng/u1	0.00
15) 4-Methylphenol-d8	8.602	113	73807	29.880	ng/u1	0.00
21) Nitrobenzene-d5	9.055	128	35724	36.458	ng/u1	0.00
24) 2-Nitrophenol-d4	9.772	143	40307	37.613	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.312	165	78503	33.638	ng/u1	0.00
31) 4-Chloroaniline-d4	10.823	131	95648	29.816	ng/u1	0.00
46) Dimethylphthalate-d6	13.932	166	297983	33.093	ng/u1	0.00
49) Acenaphthylene-d8	14.219	160	319763	31.975	ng/u1	0.00
54) 4-Nitrophenol-d4	14.725	143	44718	29.234	ng/u1	0.00
60) Fluorene-d10	15.524	176	276366	34.060	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.635	200	56266	30.016	ng/u1	0.00
73) Anthracene-d10	17.375	188	527097	34.272	ng/u1	0.00
81) Pyrene-d10	19.666	212	690180	33.449	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.372	264	724801	33.613	ng/u1	0.00

Target Compounds Qvalue

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

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