

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012987.D
 Acq On : 08 Dec 2022 16:32
 Operator : CG/JU
 Sample : PB149469BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK469

Quant Time: Dec 09 01:56:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 08 00:35:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	530216	20.000	ng/u1	0.00
20) Naphthalene-d8	10.793	136	2316472	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1593360	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	3687667	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	3514137	20.000	ng/u1	0.00
88) Perylene-d12	23.951	264	3799008	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	72136	5.817	ng/uL	0.00
4) Pyridine-d5	3.781	84	966557	27.438	ng/u1	0.00
7) Phenol-d5	7.128	99	1239199	28.694	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	805464	30.917	ng/u1	0.00
11) 2-Chlorophenol-d4	7.505	132	1014325	29.835	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	1014549	28.670	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	512068	30.087	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	577460	30.137	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.405	165	1052057	28.268	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	1492519	28.407	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	3864240	31.556	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	4042040	30.039	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	561079	25.472	ng/u1	0.00
60) Fluorene-d10	15.610	176	3281520	31.675	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	573653	24.173	ng/u1	0.00
73) Anthracene-d10	17.469	188	5499533	33.057	ng/u1	0.00
81) Pyrene-d10	19.698	212	6546329	32.454	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.786	264	6360765	33.591	ng/u1	0.00

Target Compounds Qvalue

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

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