

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013032.D
 Acq On : 10 Dec 2022 10:35
 Operator : CG/JU
 Sample : PB149532BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK532

Quant Time: Dec 11 23:48:53 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	7.969	152	591253	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.787	136	2593667	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.610	164	1740884	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	3952772	20.000	ng/u1	0.00
79) Chrysene-d12	21.451	240	3260688	20.000	ng/u1	0.00
88) Perylene-d12	23.939	264	3009795	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	70402	5.091	ng/uL	0.00
4) Pyridine-d5	3.781	84	953911	24.284	ng/u1	0.00
7) Phenol-d5	7.128	99	1267990	26.329	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	798160	27.474	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	1049473	27.682	ng/u1	-0.01
15) 4-Methylphenol-d8	8.675	113	1052280	26.667	ng/u1	-0.01
21) Nitrobenzene-d5	9.140	128	526201	27.613	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	609693	28.419	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.404	165	1096088	26.303	ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	1512913	25.717	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	3722480	27.822	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	4085654	27.790	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	564154	23.441	ng/u1	0.00
60) Fluorene-d10	15.604	176	3196202	28.237	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	564848	22.206	ng/u1	0.00
73) Anthracene-d10	17.469	188	5315807	29.809	ng/u1	0.00
81) Pyrene-d10	19.698	212	6059048	32.373	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.780	264	4592527	30.613	ng/u1	0.00

Target Compounds Qvalue

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

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