

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013123.D
 Acq On : 19 Dec 2022 12:51
 Operator : CG/JU
 Sample : PB149588BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK588

Quant Time: Dec 19 21:45:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 19 21:43:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	467132	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	1952988	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	1274265	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	2952013	20.000	ng/u1	0.00
79) Chrysene-d12	21.439	240	2813380	20.000	ng/u1	0.00
88) Perylene-d12	23.921	264	2892663	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	58213	5.328	ng/uL	0.00
4) Pyridine-d5	3.764	84	815515	26.277	ng/u1	0.00
7) Phenol-d5	7.111	99	1057655	27.797	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	697837	30.403	ng/u1	0.00
11) 2-Chlorophenol-d4	7.487	132	866114	28.916	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	843912	27.069	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	405111	28.232	ng/u1	0.00
24) 2-Nitrophenol-d4	9.852	143	457645	28.329	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	848315	27.036	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	1165133	26.303	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	2931921	29.938	ng/u1	0.00
49) Acenaphthylene-d8	14.293	160	3079955	28.621	ng/u1	0.00
54) 4-Nitrophenol-d4	14.787	143	391508	22.224	ng/u1	0.00
60) Fluorene-d10	15.592	176	2511573	30.314	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	406765	21.412	ng/u1	0.00
73) Anthracene-d10	17.451	188	4095617	30.753	ng/u1	0.00
81) Pyrene-d10	19.686	212	5005711	30.997	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.763	264	4418035	30.642	ng/u1	0.00

Target Compounds Qvalue

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

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