

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012988.D
 Acq On : 08 Dec 2022 17:06
 Operator : CG/JU
 Sample : PB149470BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK470

Quant Time: Dec 09 01:56:44 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	582889	20.000 ng/u1	0.00
20) Naphthalene-d8	10.793	136	2511487	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1726620	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.369	188	3936324	20.000 ng/u1	0.00
79) Chrysene-d12	21.457	240	3622059	20.000 ng/u1	0.00
88) Perylene-d12	23.945	264	3787639	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.358	96	83705	6.140 ng/uL	0.00
4) Pyridine-d5	3.781	84	1150810	29.717 ng/u1	0.00
7) Phenol-d5	7.128	99	1426290	30.041 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	935838	32.675 ng/u1	0.00
11) 2-Chlorophenol-d4	7.505	132	1163559	31.132 ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	1168303	30.032 ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	584475	31.674 ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	660186	31.779 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.405	165	1199211	29.720 ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	1707920	29.982 ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	4392614	33.102 ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	4576482	31.386 ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	663624	27.802 ng/u1	0.00
60) Fluorene-d10	15.610	176	3697427	32.935 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	653739	25.808 ng/u1	0.00
73) Anthracene-d10	17.469	188	6036158	33.990 ng/u1	0.00
81) Pyrene-d10	19.698	212	7088939	34.097 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.786	264	6574586	34.825 ng/u1	0.00

Target Compounds Qvalue

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

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