

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012977.D
 Acq On : 08 Dec 2022 09:16
 Operator : CG/JU
 Sample : PB149448BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK448

Quant Time: Dec 09 01:51:05 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	654754	20.000	ng/u1	0.00
20) Naphthalene-d8	10.793	136	2791015	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1903935	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.375	188	4371007	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	3936947	20.000	ng/u1	0.00
88) Perylene-d12	23.951	264	3984638	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	82804	5.407	ng/uL	0.00
4) Pyridine-d5	3.787	84	1152925	26.504	ng/u1	0.00
7) Phenol-d5	7.134	99	1378214	25.843	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	894571	27.806	ng/u1	0.00
11) 2-Chlorophenol-d4	7.505	132	1158129	27.586	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	1128768	25.831	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	565037	27.554	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	641735	27.797	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	1159857	25.866	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	1676195	26.478	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	4247940	29.031	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	4476491	27.841	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	631451	23.990	ng/u1	0.00
60) Fluorene-d10	15.610	176	3490707	28.198	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	640864	22.784	ng/u1	0.00
73) Anthracene-d10	17.469	188	5768860	29.254	ng/u1	0.00
81) Pyrene-d10	19.704	212	6798550	30.085	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.786	264	5908646	29.750	ng/u1	0.00

Target Compounds	Qvalue

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

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