

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121222\
 Data File : BP013078.D
 Acq On : 12 Dec 2022 08:38
 Operator : CG/JU
 Sample : PB149532BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK532

Quant Time: Dec 12 10:59:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	496525	20.000	ng/u1	0.00
20) Naphthalene-d8	10.787	136	2149922	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1417741	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	3252873	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	2669237	20.000	ng/u1	0.00
88) Perylene-d12	23.951	264	2643449	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	73482	6.328	ng/uL	0.00
4) Pyridine-d5	3.781	84	995374	30.174	ng/u1	0.00
7) Phenol-d5	7.128	99	1314503	32.502	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	838220	34.358	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	1074257	33.742	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	1061185	32.023	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	521498	33.014	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	596932	33.567	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.405	165	1107811	32.072	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	1497918	30.718	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	3661168	33.601	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	3959157	33.068	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	509728	26.007	ng/u1	0.00
60) Fluorene-d10	15.610	176	3155592	34.233	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	504072	24.080	ng/u1	0.00
73) Anthracene-d10	17.469	188	5240070	35.707	ng/u1	0.00
81) Pyrene-d10	19.698	212	6045472	39.458	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.786	264	4835998	36.703	ng/u1	0.01

Target Compounds Qvalue

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

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