

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012546.D
 Acq On : 08 Nov 2022 11:12
 Operator : CG/JU
 Sample : PB148826BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK826

Quant Time: Nov 08 21:40:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.769	152	333715	20.000 ng/u1	0.00
20) Naphthalene-d8	10.569	136	1593705	20.000 ng/u1	-0.01
38) Acenaphthene-d10	14.428	164	1116114	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.186	188	2544988	20.000 ng/u1	0.00
79) Chrysene-d12	21.286	240	2258548	20.000 ng/u1	0.00
88) Perylene-d12	23.662	264	2049512	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.228	96	44425	5.360 ng/uL	0.00
4) Pyridine-d5	3.652	84	599113	25.164 ng/u1	0.00
7) Phenol-d5	6.952	99	872732	29.157 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	505645	28.299 ng/u1	0.00
11) 2-Chlorophenol-d4	7.304	132	653482	29.897 ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	705483	29.801 ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	343817	33.268 ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	389598	36.320 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	702304	30.210 ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	1050162	30.339 ng/u1	0.00
46) Dimethylphthalate-d6	13.845	166	2276718	29.798 ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	2521063	28.874 ng/u1	0.00
54) 4-Nitrophenol-d4	14.657	143	313446	21.676 ng/u1	0.00
60) Fluorene-d10	15.422	176	2007909	30.695 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	254218	19.186 ng/u1	0.00
73) Anthracene-d10	17.286	188	3266069	29.784 ng/u1	0.00
81) Pyrene-d10	19.533	212	3735180	32.940 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.509	264	2974042	29.698 ng/u1	0.00

Target Compounds Qvalue

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

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