

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012556.D
 Acq On : 08 Nov 2022 18:08
 Operator : CG/JU
 Sample : PB148832BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK832

Quant Time: Nov 08 22:44:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 08 22:40:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	165366	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	823751	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	588622	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	1339610	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	1145331	20.000	ng/u1	0.00
88) Perylene-d12	23.663	264	1009817	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	32338	7.874	ng/uL	0.00
4) Pyridine-d5	3.646	84	462847	39.232	ng/u1	0.00
7) Phenol-d5	6.963	99	690877	46.580	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	371175	41.920	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	514061	47.461	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	571061	48.681	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	263528	49.333	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	298005	53.748	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	568963	47.351	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	820908	45.883	ng/u1	0.00
46) Dimethylphthalate-d6	13.840	166	1806652	44.836	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	2034779	44.188	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	267410	35.065	ng/u1	0.00
60) Fluorene-d10	15.422	176	1597050	46.293	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	276092	39.585	ng/u1	0.00
73) Anthracene-d10	17.286	188	2594778	44.954	ng/u1	0.00
81) Pyrene-d10	19.528	212	2909523	50.598	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.510	264	2255785	45.718	ng/u1	0.00

Target Compounds Qvalue

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

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