

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012263.D
 Acq On : 22 Oct 2022 01:49
 Operator : CG/JU
 Sample : PB148371BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK371

Quant Time: Oct 22 02:37:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 22 01:24:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	444114	20.000	ng/u1	0.00
20) Naphthalene-d8	10.622	136	1728173	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.469	164	950437	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1839690	20.000	ng/u1	0.00
79) Chrysene-d12	21.321	240	1886794	20.000	ng/u1	0.00
88) Perylene-d12	23.721	264	1988044	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	58818	5.360	ng/uL	0.00
4) Pyridine-d5	3.675	84	786232	24.864	ng/u1	0.00
7) Phenol-d5	6.993	99	999781	26.076	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	617080	26.538	ng/u1	0.00
11) 2-Chlorophenol-d4	7.351	132	810891	27.785	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	749979	25.168	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	382675	29.915	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	404749	29.326	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	682676	26.370	ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	995996	27.078	ng/u1	0.00
46) Dimethylphthalate-d6	13.886	166	1857272	28.092	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	2155266	28.560	ng/u1	0.00
54) 4-Nitrophenol-d4	14.686	143	290980	23.492	ng/u1	0.00
60) Fluorene-d10	15.469	176	1568858	27.583	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	153918	14.707	ng/u1	0.00
73) Anthracene-d10	17.327	188	2399230	29.753	ng/u1	0.00
81) Pyrene-d10	19.569	212	2813918	26.194	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.568	264	3048945	30.582	ng/u1	0.00

Target Compounds Qvalue

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

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