

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012428.D
 Acq On : 01 Nov 2022 15:34
 Operator : CG/JU
 Sample : PB148597BL
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK597

Quant Time: Nov 01 22:59:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	298377	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1336889	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.440	164	887357	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	1939729	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	1897084	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	1636130	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	51550	6.956	ng/uL	0.00
4) Pyridine-d5	3.658	84	670551	31.500	ng/u1	0.00
7) Phenol-d5	6.963	99	906197	33.861	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	561238	35.130	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	698820	35.757	ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	726634	34.330	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	373219	43.050	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	406519	45.177	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	677744	34.754	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	1027557	35.389	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2314342	38.100	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2611051	37.614	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	405617	35.282	ng/u1	0.00
60) Fluorene-d10	15.440	176	2000916	38.473	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	345088	34.170	ng/u1	0.00
73) Anthracene-d10	17.304	188	3210542	38.413	ng/u1	0.00
81) Pyrene-d10	19.545	212	3770286	39.585	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3173694	39.699	ng/u1	-0.01

Target Compounds Qvalue

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

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