

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012246.D
 Acq On : 21 Oct 2022 15:31
 Operator : CG/JU
 Sample : PB148371BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK371

Quant Time: Oct 21 22:32:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	388416	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1652153	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	944259	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	1663457	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1326485	20.000	ng/u1	0.00
88) Perylene-d12	23.727	264	1446742	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	51423	5.358	ng/uL	0.00
4) Pyridine-d5	3.681	84	715831	25.884	ng/u1	0.00
7) Phenol-d5	6.993	99	976739	29.128	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	575373	28.293	ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	777614	30.466	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	769167	29.513	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	394965	32.296	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	439144	33.283	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	727493	29.394	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	1059384	30.127	ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	1963615	29.895	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2317565	30.911	ng/u1	0.00
54) 4-Nitrophenol-d4	14.693	143	274600	22.315	ng/u1	0.00
60) Fluorene-d10	15.469	176	1658573	29.351	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	182082	19.241	ng/u1	0.00
73) Anthracene-d10	17.334	188	2302630	31.580	ng/u1	0.00
81) Pyrene-d10	19.575	212	2272032	30.084	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	2319496	31.970	ng/u1	-0.01

Target Compounds	Qvalue

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

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