

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012444.D
 Acq On : 02 Nov 2022 00:45
 Operator : CG/JU
 Sample : PB148613BL
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK613

Quant Time: Nov 02 01:54:41 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	7.787	152	267778	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1193792	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.440	164	815244	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1812546	20.000	ng/u1	-0.01
79) Chrysene-d12	21.298	240	1673307	20.000	ng/u1	0.00
88) Perylene-d12	23.680	264	1420715	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	43446	6.533	ng/uL	0.00
4) Pyridine-d5	3.658	84	561595	29.397	ng/u1	0.00
7) Phenol-d5	6.964	99	799716	33.297	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	465868	32.492	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	618996	35.292	ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	649181	34.175	ng/u1	0.00
21) Nitrobenzene-d5	8.958	128	329866	42.610	ng/u1	0.00
24) 2-Nitrophenol-d4	9.675	143	376679	46.879	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.216	165	640311	36.771	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	892736	34.431	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2074439	37.171	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2317727	36.342	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	363282	34.394	ng/u1	0.00
60) Fluorene-d10	15.440	176	1785651	37.371	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	187451	19.864	ng/u1	0.00
73) Anthracene-d10	17.298	188	2895483	37.075	ng/u1	-0.01
81) Pyrene-d10	19.539	212	3344101	39.805	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.527	264	2643506	38.081	ng/u1	-0.02

Target Compounds Qvalue

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

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