

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102523\
 Data File : BP017809.D
 Acq On : 25 Oct 2023 18:42
 Operator : MA/JU
 Sample : PB156616BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK616

Quant Time: Oct 26 05:30:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	138942	20.000	ng/u1	0.00
20) Naphthalene-d8	10.710	136	508093	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.581	164	285277	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	585191	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.510	240	510292	20.000	ng/u1	0.00
88) Perylene-d12	24.062	264	519710	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	8491	2.273	ng/uL	0.02
4) Pyridine-d5	3.711	84	101007	10.329	ng/u1	0.02
7) Phenol-d5	7.046	99	121525	10.511	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	77999	10.924	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	99913	10.595	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	130913	14.418	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	55080	13.378	ng/u1	0.00
24) 2-Nitrophenol-d4	9.804	143	66726	14.668	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	137748	15.437	ng/u1	0.00
31) 4-Chloroaniline-d4	10.887	131	264938	22.687	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	926258	38.678	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	858669	32.151	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	124578	36.508	ng/u1	0.00
60) Fluorene-d10	15.581	176	719358	34.749	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	137370	35.638	ng/u1	0.00
73) Anthracene-d10	17.469	188	1233297	41.241	ng/u1	0.00
81) Pyrene-d10	19.727	212	1412756	44.159	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	1253605	43.308	ng/u1	0.00

Target Compounds Qvalue

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

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