

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017774.D
 Acq On : 24 Oct 2023 10:38
 Operator : MA/JU
 Sample : PB156501BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK501

Quant Time: Oct 25 01:49:27 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	7.892	152	145174	20.000	ng/u1	0.00
20) Naphthalene-d8	10.728	136	589252	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.592	164	361082	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	796807	20.000	ng/u1	0.01
79) Chrysene-d12	21.527	240	635815	20.000	ng/u1	0.01
88) Perylene-d12	24.092	264	633827	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	25762	6.602	ng/uL	0.00
4) Pyridine-d5	3.693	84	327026	32.008	ng/u1	0.00
7) Phenol-d5	7.051	99	381572	31.585	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	252041	33.782	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	311311	31.595	ng/u1	0.00
15) 4-Methylphenol-d8	8.616	113	296897	31.296	ng/u1	0.00
21) Nitrobenzene-d5	9.104	128	146390	30.659	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	166136	31.491	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	285992	27.635	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	364325	26.901	ng/u1	0.01
46) Dimethylphthalate-d6	14.010	166	892716	29.451	ng/u1	0.00
49) Acenaphthylene-d8	14.292	160	1014051	29.998	ng/u1	0.01
54) 4-Nitrophenol-d4	14.839	143	102119	23.644	ng/u1	0.01
60) Fluorene-d10	15.598	176	815337	31.117	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	113194	21.567	ng/u1	0.01
73) Anthracene-d10	17.486	188	1215728	29.857	ng/u1	0.01
81) Pyrene-d10	19.745	212	1345244	33.747	ng/u1	0.01
92) Benzo(a)pyrene-d12	23.921	264	1079094	30.567	ng/u1	0.02

Target Compounds Qvalue

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

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