

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101023\
 Data File : BP017631.D
 Acq On : 10 Oct 2023 11:01
 Operator : MA/JU
 Sample : PB156093BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK093

Quant Time: Oct 10 23:57:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 04:45:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	87551	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	364701	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.622	164	228772	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	549922	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	552448	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	613397	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	19532	9.051	ng/uL	0.00
4) Pyridine-d5	3.716	84	256388	42.205	ng/u1	0.00
7) Phenol-d5	7.081	99	298619	40.332	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	191930	41.799	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	249331	42.811	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	237526	39.828	ng/u1	0.00
21) Nitrobenzene-d5	9.134	128	115575	41.127	ng/u1	0.00
24) 2-Nitrophenol-d4	9.851	143	129056	41.814	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	226281	37.749	ng/u1	0.00
31) 4-Chloroaniline-d4	10.934	131	336580	39.423	ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	819389	45.111	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	844350	42.696	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	94011	34.149	ng/u1	0.00
60) Fluorene-d10	15.628	176	704612	44.357	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	105137	32.232	ng/u1	0.00
73) Anthracene-d10	17.516	188	1200620	45.493	ng/u1	0.00
81) Pyrene-d10	19.769	212	1478137	46.907	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.957	264	1461489	45.957	ng/u1	0.00

Target Compounds	Qvalue

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

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