

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101723\
 Data File : BP017703.D
 Acq On : 17 Oct 2023 18:07
 Operator : MA/JU
 Sample : PB156373BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK373

Quant Time: Oct 18 01:01:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 16:14:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	124399	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	495839	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.604	164	326138	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	798673	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	904392	20.000	ng/u1	0.00
88) Perylene-d12	24.110	264	1071738	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	28120	8.334	ng/uL	0.00
4) Pyridine-d5	3.699	84	388244	41.625	ng/u1	0.00
7) Phenol-d5	7.063	99	483914	41.783	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	302982	43.138	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	409315	45.659	ng/u1	0.00
15) 4-Methylphenol-d8	8.628	113	389639	42.120	ng/u1	0.00
21) Nitrobenzene-d5	9.116	128	195160	46.860	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	216835	46.297	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	396915	45.091	ng/u1	0.00
31) 4-Chloroaniline-d4	10.916	131	563750	44.681	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	1363667	48.278	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	1443154	47.446	ng/u1	0.00
54) 4-Nitrophenol-d4	14.857	143	188227	39.845	ng/u1	0.00
60) Fluorene-d10	15.610	176	1206047	50.196	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	196337	34.780	ng/u1	0.00
73) Anthracene-d10	17.498	188	2005400	48.639	ng/u1	0.00
81) Pyrene-d10	19.757	212	2566860	47.157	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.939	264	2986154	49.632	ng/u1	0.00

Target Compounds Qvalue

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

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