

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
 Data File : BP017923.D
 Acq On : 07 Nov 2023 09:20
 Operator : MA/JU
 Sample : PB156841BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK841

Quant Time: Nov 07 21:54:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 07 03:40:56 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	72971	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	292707	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	187566	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	425099	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	389088	20.000	ng/u1	0.00
88) Perylene-d12	23.892	264	415543	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	9061	4.752	ng/uL	0.00
4) Pyridine-d5	3.681	84	116742	22.063	ng/u1	0.00
7) Phenol-d5	7.028	99	134950	19.721	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	90286	21.873	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	115974	22.056	ng/u1	0.00
15) 4-Methylphenol-d8	8.581	113	109770	19.408	ng/u1	0.00
21) Nitrobenzene-d5	9.063	128	54384	23.249	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	62605	23.237	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	101663	20.583	ng/u1	0.00
31) 4-Chloroaniline-d4	10.851	131	143598	20.063	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	397507	24.169	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	408712	22.647	ng/u1	0.00
54) 4-Nitrophenol-d4	14.781	143	36225	14.325	ng/u1	0.00
60) Fluorene-d10	15.522	176	323705	23.515	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	55199	17.756	ng/u1	0.00
73) Anthracene-d10	17.398	188	515513	23.324	ng/u1	0.00
81) Pyrene-d10	19.645	212	595746	24.120	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.727	264	556929	23.986	ng/u1	0.00

Target Compounds Qvalue

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

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