

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
 Data File : BP018639.D
 Acq On : 06 Dec 2023 16:21
 Operator : MA/JU
 Sample : PB157358BL
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK358

Quant Time: Dec 06 21:42:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	405585	20.000	ng/ul	0.00
20) Naphthalene-d8	10.646	136	1926124	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.481	164	1139579	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	2417549	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1813143	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	2063964	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	63489	5.334	ng/uL	0.00
4) Pyridine-d5	3.681	84	813948	25.532	ng/ul	0.00
7) Phenol-d5	7.022	99	1137977	27.959	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	671692	27.844	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	878151	27.650	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	924106	28.055	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	452530	26.397	ng/ul	0.00
24) 2-Nitrophenol-d4	9.734	143	497786	28.007	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	912194	25.100	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	994074	20.930	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	2905795	28.570	ng/ul	0.00
49) Acenaphthylene-d8	14.175	160	2823910	25.215	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	384237	23.015	ng/ul	0.00
60) Fluorene-d10	15.475	176	2343548	27.450	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	302788	20.888	ng/ul	0.00
73) Anthracene-d10	17.328	188	3314019	26.075	ng/ul	0.00
81) Pyrene-d10	19.563	212	3318918	23.515	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	3102845	25.783	ng/ul	0.00

Target Compounds	Qvalue

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

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