

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018396.D
 Acq On : 27 Nov 2023 12:41
 Operator : MA/JU
 Sample : PB157279BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK279

Quant Time: Nov 27 22:14:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 27 21:50:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	324347	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.657	136	1186807	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	609056	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1215499	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	1256768	20.000	ng/u1	0.00
88) Perylene-d12	23.880	264	1451045	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	57707	6.664	ng/uL	0.00
4) Pyridine-d5	3.675	84	700830	29.449	ng/u1	-0.01
7) Phenol-d5	7.016	99	819467	27.645	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.187	67	496657	28.020	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.381	132	718078	28.853	ng/u1	-0.01
15) 4-Methylphenol-d8	8.563	113	624605	26.110	ng/u1	-0.01
21) Nitrobenzene-d5	9.016	128	310835	32.245	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	284704	31.980	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	602653	28.059	ng/u1	0.00
31) 4-Chloroaniline-d4	10.793	131	846233	29.213	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1612580	29.756	ng/u1	0.00
49) Acenaphthylene-d8	14.187	160	1871573	30.308	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	201128	26.883	ng/u1	0.00
60) Fluorene-d10	15.492	176	1383785	30.528	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	122789	23.730	ng/u1	0.00
73) Anthracene-d10	17.351	188	2121557	32.656	ng/u1	0.00
81) Pyrene-d10	19.598	212	2605556	26.049	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.715	264	2773284	32.966	ng/u1	0.02

Target Compounds Qvalue

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

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