

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018317.D
 Acq On : 23 Nov 2023 15:33
 Operator : MA/JU
 Sample : PB157173BL
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK173

Quant Time: Nov 23 23:52:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	456539	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.675	136	1938266	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.504	164	1284952	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	2638327	20.000	ng/ul	-0.01
79) Chrysene-d12	21.357	240	1649781	20.000	ng/ul	0.00
88) Perylene-d12	23.780	264	1647198	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	82361	6.757	ng/uL	0.00
4) Pyridine-d5	3.687	84	1141290	34.071	ng/ul	0.00
7) Phenol-d5	7.034	99	1402909	33.624	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	855031	34.270	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	1149155	32.804	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	1161792	34.504	ng/ul	-0.01
21) Nitrobenzene-d5	9.034	128	573050	36.400	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	561801	38.640	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	1153235	32.876	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.810	131	1759351	37.188	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	3941696	34.475	ng/ul	-0.01
49) Acenaphthylene-d8	14.198	160	4415185	33.890	ng/ul	0.00
54) 4-Nitrophenol-d4	14.693	143	496926	31.483	ng/ul	0.00
60) Fluorene-d10	15.498	176	3260359	34.093	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	323832	28.833	ng/ul	0.00
73) Anthracene-d10	17.351	188	4779722	33.895	ng/ul	-0.01
81) Pyrene-d10	19.592	212	4790623	36.485	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.622	264	3332150	34.893	ng/ul	0.00

Target Compounds Qvalue

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

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