

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019253.D
 Acq On : 03 Jan 2024 17:50
 Operator : MA/JU
 Sample : PB158128BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK128

Quant Time: Jan 04 04:49:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	175647	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	631848	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.439	164	384562	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	922214	20.000	ng/u1	0.00
79) Chrysene-d12	21.304	240	1043237	20.000	ng/u1	0.00
88) Perylene-d12	23.668	264	1220612	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	28204	5.484	ng/uL	0.00
4) Pyridine-d5	3.658	84	380190	29.453	ng/u1	0.00
7) Phenol-d5	6.987	99	433642	27.530	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	243222	26.691	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	343258	27.031	ng/u1	0.00
15) 4-Methylphenol-d8	8.528	113	321839	26.244	ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	153064	28.024	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	173839	29.449	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	325678	27.322	ng/u1	0.00
31) 4-Chloroaniline-d4	10.745	131	414809	28.292	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	888037	26.144	ng/u1	-0.01
49) Acenaphthylene-d8	14.134	160	1048242	28.441	ng/u1	0.00
54) 4-Nitrophenol-d4	14.681	143	156099	27.215	ng/u1	0.02
60) Fluorene-d10	15.439	176	813867	27.734	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	92696	15.332	ng/u1	0.00
73) Anthracene-d10	17.298	188	1367176	28.388	ng/u1	0.00
81) Pyrene-d10	19.545	212	1760688	27.693	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.515	264	1933405	27.557	ng/u1	0.00

Target Compounds Qvalue

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

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