

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091224\
 Data File : BP021892.D
 Acq On : 12 Sep 2024 11:56
 Operator : RC/JU
 Sample : PB163193BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK193

Quant Time: Sep 13 02:36:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 12 12:48:41 2024
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.763	152	107735	20.000 ng/u1	0.03
20) Naphthalene-d8	10.528	136	373354	20.000 ng/u1	0.02
38) Acenaphthene-d10	14.363	164	308558	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.157	188	746377	20.000 ng/u1	# 0.00
79) Chrysene-d12	21.580	240	842360	20.000 ng/u1	#-0.01
88) Perylene-d12	24.904	264	964123	20.000 ng/u1	-0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.269	96	13710	4.116 ng/uL	0.01
4) Pyridine-d5	3.681	84	179297	18.285 ng/u1	0.02
7) Phenol-d5	6.934	99	217463	23.623 ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.099	67	126322	23.626 ng/u1	0.03
11) 2-Chlorophenol-d4	7.299	132	183142	28.580 ng/u1	0.02
15) 4-Methylphenol-d8	8.457	113	164643	22.401 ng/u1	0.03
21) Nitrobenzene-d5	8.899	128	81360	34.990 ng/u1	0.02
24) 2-Nitrophenol-d4	9.616	143	90474	34.227 ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.151	165	170862	29.189 ng/u1	0.02
31) 4-Chloroaniline-d4	10.657	131	221391	25.951 ng/u1	0.02
46) Dimethylphthalate-d6	13.775	166	730936	31.531 ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	722329	28.406 ng/u1	0.00
54) 4-Nitrophenol-d4	14.563	143	112242	25.613 ng/u1	0.00
60) Fluorene-d10	15.369	176	617073	28.005 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	112833	26.424 ng/u1	0.00
73) Anthracene-d10	17.257	188	1019326	29.400 ng/u1	0.00
81) Pyrene-d10	19.622	212	1378442	29.577 ng/u1	0.00
92) Benzo(a)pyrene-d12	24.680	264	1497031	29.836 ng/u1	-0.01

Target Compounds Qvalue

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

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