

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
 Data File : BP021820.D
 Acq On : 04 Sep 2024 17:04
 Operator : RC/JU
 Sample : PB163094BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK094

Quant Time: Sep 05 05:09:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 05 05:03:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	340907	20.000	ng/u1	0.00
20) Naphthalene-d8	10.552	136	1416481	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.410	164	951340	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.210	188	2125801	20.000	ng/u1	0.00
79) Chrysene-d12	21.663	240	2033647	20.000	ng/u1	-0.02
88) Perylene-d12	25.080	264	2190495	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	57976	5.646	ng/uL	0.00
4) Pyridine-d5	3.681	84	764396	25.737	ng/u1	0.00
7) Phenol-d5	6.946	99	956253	26.284	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.105	67	538002	27.173	ng/u1	0.01
11) 2-Chlorophenol-d4	7.311	132	633545	26.973	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	723225	25.725	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	320386	28.114	ng/u1	0.00
24) 2-Nitrophenol-d4	9.634	143	338993	29.305	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	594220	26.067	ng/u1	0.00
31) 4-Chloroaniline-d4	10.687	131	848591	25.587	ng/u1	0.00
46) Dimethylphthalate-d6	13.816	166	2158222	29.648	ng/u1	0.00
49) Acenaphthylene-d8	14.104	160	2272595	27.807	ng/u1	0.01
54) 4-Nitrophenol-d4	14.616	143	373965	27.003	ng/u1	0.00
60) Fluorene-d10	15.422	176	1781482	29.110	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.534	200	301524	25.789	ng/u1	0.00
73) Anthracene-d10	17.310	188	2853643	28.285	ng/u1	0.00
81) Pyrene-d10	19.686	212	3475083	29.893	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.833	264	3322639	28.446	ng/u1	-0.02

Target Compounds Qvalue

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

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