

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
 Data File : BP020152.D
 Acq On : 02 May 2024 19:12
 Operator : MA/JU
 Sample : PB160609BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: May 03 00:42:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 03 00:38:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	91592	20.000	ng/ul	0.00
20) Naphthalene-d8	10.399	136	387802	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	280210	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.140	188	647510	20.000	ng/ul	0.00
79) Chrysene-d12	21.639	240	580622	20.000	ng/ul	0.00
88) Perylene-d12	25.010	264	542068	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	13711	6.322	ng/uL	0.00
4) Pyridine-d5	3.628	84	63244	10.036	ng/ul	0.00
7) Phenol-d5	6.811	99	240109	30.526	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	154513	32.337	ng/ul	0.00
11) 2-Chlorophenol-d4	7.164	132	181213	32.227	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113	198651	31.238	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	96664	33.287	ng/ul	0.00
24) 2-Nitrophenol-d4	9.499	143	111891	33.646	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	181236	29.865	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	40	0.005	ng/ul	-0.02
46) Dimethylphthalate-d6	13.722	166	746636	36.488	ng/ul	0.00
49) Acenaphthylene-d8	13.993	160	770202	34.073	ng/ul	0.00
54) 4-Nitrophenol-d4	14.563	143	93540	29.743	ng/ul	0.02
60) Fluorene-d10	15.334	176	612151	35.866	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.487	200	131090	31.904	ng/ul	0.01
73) Anthracene-d10	17.251	188	1022492	36.502	ng/ul	0.01
81) Pyrene-d10	19.651	212	1219473	35.930	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.774	264	958519	36.571	ng/ul	0.00

Target Compounds Qvalue

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

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