

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046569.D
 Acq On : 13 Jul 2024 05:57
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4165

Quant Time: Jul 13 07:04:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 05:09:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	3179	0.400	ng/ul	0.00
4) Naphthalene-d8	10.266	136	8154	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.151	164	4943	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.901	188	10230	0.400	ng/ul	0.00
17) Chrysene-d12	21.105	240	7582	0.400	ng/ul	0.00
23) Perylene-d12	23.241	264	8652	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.118	96	952	0.364	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.866	152	5104	0.388	ng/ul	0.00
18) Fluoranthene-d10	18.937	212	9750	0.423	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.152	88	1034	0.371	ng/ul#	75
5) Naphthalene	10.315	128	8390	0.404	ng/ul	97
7) 2-Methylnaphthalene	11.943	142	5749	0.394	ng/ul	100
8) 1-Methylnaphthalene	12.163	142	6117	0.411	ng/ul	98
10) Acenaphthylene	13.864	152	9550	0.424	ng/ul	98
11) Acenaphthene	14.211	153	6415	0.426	ng/ul	96
12) Fluorene	15.206	166	7364	0.395	ng/ul	97
14) Pentachlorophenol	16.555	266	1507	0.291	ng/ul	98
15) Phenanthrene	16.943	178	11698	0.401	ng/ul	99
16) Anthracene	17.032	178	11205	0.393	ng/ul	98
19) Fluoranthene	18.970	202	13772	0.442	ng/ul	99
20) Pyrene	19.332	202	14062	0.417	ng/ul	99
21) Benzo(a)anthracene	21.088	228	13160	0.392	ng/ul	99
22) Chrysene	21.140	228	12562	0.387	ng/ul	100
24) Benzo(b)fluoranthene	22.610	252	13557	0.392	ng/ul	97
25) Benzo(k)fluoranthene	22.651	252	13283	0.391	ng/ul	96
26) Benzo(a)pyrene	23.151	252	12562	0.442	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.360	276	16425	0.379	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.383	278	13032	0.380	ng/ul	98
29) Benzo(g,h,i)perylene	26.004	276	13525	0.392	ng/ul	99

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

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