

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011223\
 Data File : BM038432.D
 Acq On : 12 Jan 2023 09:28
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4063

Quant Time: Jan 12 10:29:28 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM123022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 12 01:15:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.996	152	1449	0.400	ng/u1	0.00
4) Naphthalene-d8	10.806	136	4104	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.629	164	2266	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.371	188	4830	0.400	ng/u1	0.00
17) Chrysene-d12	21.544	240	5092	0.400	ng/u1	0.00
23) Perylene-d12	23.974	264	5717	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.368	96	864	0.400	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.390	152	2153	0.350	ng/u1	0.00
18) Fluoranthene-d10	19.389	212	5137	0.356	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.402	88	871	0.393	ng/u1	98
5) Naphthalene	10.856	128	4011	0.366	ng/u1	98
7) 2-Methylnaphthalene	12.467	142	2377	0.353	ng/u1	98
8) 1-Methylnaphthalene	12.682	142	2473	0.361	ng/u1	98
10) Acenaphthylene	14.351	152	3783	0.359	ng/u1	99
11) Acenaphthene	14.689	153	2842	0.363	ng/u1	98
12) Fluorene	15.674	166	3176	0.358	ng/u1	95
14) Pentachlorophenol	17.029	266	634	0.368	ng/u1	97
15) Phenanthrene	17.413	178	5006	0.351	ng/u1	99
16) Anthracene	17.506	178	4609	0.351	ng/u1	98
19) Fluoranthene	19.422	202	6422	0.365	ng/u1	99
20) Pyrene	19.784	202	6789	0.361	ng/u1	99
21) Benzo(a)anthracene	21.527	228	6386	0.377	ng/u1	99
22) Chrysene	21.580	228	6652	0.372	ng/u1	98
24) Benzo(b)fluoranthene	23.225	252	7789	0.376	ng/u1	95
25) Benzo(k)fluoranthene	23.278	252	7951	0.385	ng/u1	95
26) Benzo(a)pyrene	23.862	252	7496	0.383	ng/u1	93
27) Indeno(1,2,3-cd)pyrene	26.505	276	10149	0.355	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.515	278	8072	0.360	ng/u1	96
29) Benzo(g,h,i)perylene	27.283	276	8927	0.351	ng/u1	99

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

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