

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032093.D
 Acq On : 05 Sep 2021 01:14
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4123

Quant Time: Sep 06 04:26:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM090321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 16:21:27 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.452	152	1578	0.400	ng/ul	0.00
4) Naphthalene-d8	10.208	136	4435	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.095	164	1592	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.855	188	3135	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.059	240	2458	0.400	ng/ul	#-0.01
23) Perylene-d12	23.173	264	2998	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.068	96	740	0.427	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.809	152	2001	0.292	ng/ul	-0.01
18) Fluoranthene-d10	18.892	212	2800	0.389	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.099	88	728	0.430	ng/ul#	89
5) Naphthalene	10.257	128	4664	0.396	ng/ul	99
7) 2-Methylnaphthalene	11.881	142	2479	0.302	ng/ul	97
8) 1-Methylnaphthalene	12.101	142	2344	0.290	ng/ul	94
10) Acenaphthylene	13.815	152	3355	0.522	ng/ul	100
11) Acenaphthene	14.159	153	2456	0.467	ng/ul#	84
12) Fluorene	15.160	166	2629	0.429	ng/ul	93
14) Pentachlorophenol	16.522	266	571	0.689	ng/ul	96
15) Phenanthrene	16.897	178	3916	0.410	ng/ul	99
16) Anthracene	16.990	178	3477	0.415	ng/ul	98
19) Fluoranthene	18.923	202	4177	0.393	ng/ul#	96
20) Pyrene	19.289	202	4127	0.377	ng/ul#	95
21) Benzo(a)anthracene	21.045	228	4023	0.448	ng/ul	100
22) Chrysene	21.098	228	4282	0.415	ng/ul	98
24) Benzo(b)fluoranthene	22.548	252	6162	0.447	ng/ul	95
25) Benzo(k)fluoranthene	22.589	252	5945	0.418	ng/ul	96
26) Benzo(a)pyrene	23.084	252	5904	0.490	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	25.245	276	7677	0.468	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.252	278	5851	0.444	ng/ul	95
29) Benzo(g,h,i)perylene	25.872	276	6728	0.461	ng/ul	96

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

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