

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101821\
 Data File : BM032564.D
 Acq On : 18 Oct 2021 09:23
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4141

Quant Time: Oct 18 10:11:19 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 15 12:37:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.553	152	4800	0.40	ng/ul	0.00
4) Naphthalene-d8	10.335	136	15036	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.209	164	7585	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.946	188	14811	0.40	ng/ul	# 0.00
17) Chrysene-d12	21.123	240	9968	0.40	ng/ul	-0.01
23) Perylene-d12	23.259	264	7962	0.40	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.924	96	2242	0.42	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.931	152	7659	0.38	ng/ul	0.00
18) Fluoranthene-d10	18.970	212	13621	0.43	ng/ul	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.958	88	2573	0.43	ng/ul#	81
5) Naphthalene	10.384	128	18059	0.40	ng/ul	99
7) 2-Methylnaphthalene	12.008	142	11533	0.39	ng/ul	99
8) 1-Methylnaphthalene	12.228	142	11411	0.39	ng/ul	100
10) Acenaphthylene	13.928	152	12286	0.38	ng/ul	99
11) Acenaphthene	14.270	153	12152	0.41	ng/ul	99
12) Fluorene	15.256	166	13384	0.40	ng/ul	99
14) Pentachlorophenol	16.616	266	1751	0.46	ng/ul	99
15) Phenanthrene	16.984	178	20587	0.41	ng/ul	99
16) Anthracene	17.075	178	16682	0.38	ng/ul	99
19) Fluoranthene	19.000	202	21818	0.44	ng/ul	100
20) Pyrene	19.362	202	21642	0.43	ng/ul	100
21) Benzo(a)anthracene	21.109	228	15017	0.39	ng/ul	99
22) Chrysene	21.159	228	17994	0.42	ng/ul	99
24) Benzo(b)fluoranthene	22.627	252	16084	0.44	ng/ul	97
25) Benzo(k)fluoranthene	22.665	252	16487	0.45	ng/ul	97
26) Benzo(a)pyrene	23.167	252	12777	0.41	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.360	276	16377	0.43	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.364	278	12948	0.44	ng/ul	99
29) Benzo(g,h,i)perylene	26.002	276	14760	0.45	ng/ul	99

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

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