

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100821\
 Data File : BM032392.D
 Acq On : 08 Oct 2021 14:11
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4136

Quant Time: Oct 08 14:48:36 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM092721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 08 12:27:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.619	152	7419	0.40	ng/ul	0.00
4) Naphthalene-d8	10.406	136	25374	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.266	164	13476	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.995	188	27077	0.40	ng/ul	0.00
17) Chrysene-d12	21.157	240	23226	0.40	ng/ul	0.00
23) Perylene-d12	23.300	264	20174	0.40	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.965	96	3481	0.44	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.999	152	12987	0.38	ng/ul	0.00
18) Fluoranthene-d10	19.015	212	25348	0.40	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.999	88	3759	0.46	ng/ul	92
5) Naphthalene	10.456	128	28455	0.38	ng/ul	100
7) 2-Methylnaphthalene	12.075	142	18636	0.39	ng/ul	100
8) 1-Methylnaphthalene	12.295	142	18473	0.39	ng/ul	100
10) Acenaphthylene	13.987	152	18245	0.34	ng/ul#	87
11) Acenaphthene	14.327	153	19546	0.39	ng/ul	98
12) Fluorene	15.309	166	21614	0.38	ng/ul	99
14) Pentachlorophenol	16.661	266	2730	0.40	ng/ul	99
15) Phenanthrene	17.033	178	34174	0.39	ng/ul	100
16) Anthracene	17.123	178	27734	0.38	ng/ul	100
19) Fluoranthene	19.042	202	38451	0.40	ng/ul	97
20) Pyrene	19.404	202	39144	0.40	ng/ul	100
21) Benzo(a)anthracene	21.142	228	31633	0.39	ng/ul	100
22) Chrysene	21.191	228	36759	0.38	ng/ul	100
24) Benzo(b)fluoranthene	22.663	252	37276	0.36	ng/ul	96
25) Benzo(k)fluoranthene	22.702	252	37410	0.37	ng/ul	96
26) Benzo(a)pyrene	23.208	252	29922	0.38	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.418	276	36518	0.36	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.420	278	29616	0.36	ng/ul	97
29) Benzo(g,h,i)perylene	26.069	276	32016	0.34	ng/ul	99

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

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