

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
 Data File : BM031652.D
 Acq On : 10 Aug 2021 22:23
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4110

Quant Time: Aug 11 05:58:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 10 18:36:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.576	152	1643	0.400	ng/ul	0.00
4) Naphthalene-d8	10.338	136	6139	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.212	164	3277	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.964	188	6092	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.156	240	4985	0.400	ng/ul	# 0.00
23) Perylene-d12	23.314	264	5040	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	681	0.373	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.935	152	3709	0.359	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	6399	0.385	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.220	88	662	0.359	ng/ul#	74
5) Naphthalene	10.388	128	6883	0.376	ng/ul	99
7) 2-Methylnaphthalene	12.007	142	4528	0.362	ng/ul	99
8) 1-Methylnaphthalene	12.232	142	4358	0.353	ng/ul	99
10) Acenaphthylene	13.932	152	5683	0.407	ng/ul	100
11) Acenaphthene	14.273	153	4501	0.383	ng/ul	99
12) Fluorene	15.270	166	4854	0.359	ng/ul	99
14) Pentachlorophenol	16.624	266	746	0.382	ng/ul	96
15) Phenanthrene	17.006	178	7285	0.377	ng/ul	100
16) Anthracene	17.100	178	6974	0.385	ng/ul	99
19) Fluoranthene	19.028	202	8151	0.383	ng/ul	96
20) Pyrene	19.390	202	8347	0.387	ng/ul#	93
21) Benzo(a)anthracene	21.142	228	7617	0.378	ng/ul	98
22) Chrysene	21.193	228	7767	0.373	ng/ul	99
24) Benzo(b)fluoranthene	22.672	252	7894	0.348	ng/ul	91
25) Benzo(k)fluoranthene	22.718	252	7825	0.334	ng/ul#	94
26) Benzo(a)pyrene	23.224	252	7113	0.358	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.453	276	8984	0.349	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.468	278	7216	0.348	ng/ul#	95
29) Benzo(g,h,i)perylene	26.100	276	7480	0.341	ng/ul#	92

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

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