

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033879.D
 Acq On : 26 Jan 2022 20:57
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4055

Quant Time: Jan 27 02:13:38 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.939	152	1788	0.400	ng/ul	0.00
4) Naphthalene-d8	10.752	136	6048	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.576	164	2897	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	6003	0.400	ng/ul	0.00
17) Chrysene-d12	21.475	240	5225	0.400	ng/ul	# 0.00
23) Perylene-d12	23.873	264	5229	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	821	0.439	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	3261	0.391	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	6474	0.402	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	832	0.425	ng/ul#	79
5) Naphthalene	10.801	128	6943	0.418	ng/ul#	90
7) 2-Methylnaphthalene	12.416	142	4271	0.395	ng/ul	98
8) 1-Methylnaphthalene	12.632	142	4215	0.398	ng/ul	98
10) Acenaphthylene	14.299	152	5204	0.402	ng/ul#	94
11) Acenaphthene	14.640	153	4246	0.410	ng/ul	95
12) Fluorene	15.626	166	4789	0.405	ng/ul	98
14) Pentachlorophenol	16.973	266	997	0.361	ng/ul	98
15) Phenanthrene	17.358	178	7837	0.410	ng/ul	95
16) Anthracene	17.448	178	7030	0.393	ng/ul	94
19) Fluoranthene	19.365	202	9512	0.399	ng/ul	99
20) Pyrene	19.723	202	9900	0.406	ng/ul	95
21) Benzo(a)anthracene	21.463	228	8249	0.386	ng/ul	98
22) Chrysene	21.514	228	9235	0.408	ng/ul	98
24) Benzo(b)fluoranthene	23.141	252	9497	0.404	ng/ul	91
25) Benzo(k)fluoranthene	23.187	252	9816	0.417	ng/ul#	93
26) Benzo(a)pyrene	23.769	252	8389	0.412	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.378	276	11227	0.414	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.398	278	8690	0.407	ng/ul	94
29) Benzo(g,h,i)perylene	27.144	276	9728	0.423	ng/ul	93

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

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