

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033197.D
 Acq On : 20 Nov 2021 04:27
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4179

Quant Time: Nov 20 05:11:58 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 15:41:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.960	152	3086	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	8586	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.579	164	5236	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.312	188	11549	0.400	ng/ul	0.00
17) Chrysene-d12	21.467	240	9004	0.400	ng/ul	0.00
23) Perylene-d12	23.809	264	7329	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.417	96	1113	0.384	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	4879	0.410	ng/ul	0.00
18) Fluoranthene-d10	19.330	212	11620	0.445	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.455	88	1237	0.379	ng/ul	95
5) Naphthalene	10.805	128	10659	0.401	ng/ul	98
7) 2-Methylnaphthalene	12.411	142	6999	0.407	ng/ul	99
8) 1-Methylnaphthalene	12.627	142	7068	0.415	ng/ul	100
10) Acenaphthylene	14.299	152	10093	0.412	ng/ul#	99
11) Acenaphthene	14.640	153	8098	0.403	ng/ul	100
12) Fluorene	15.622	166	9327	0.412	ng/ul	98
14) Pentachlorophenol	16.958	266	1350	0.350	ng/ul	97
15) Phenanthrene	17.354	178	14940	0.357	ng/ul	99
16) Anthracene	17.448	178	13363	0.395	ng/ul	99
19) Fluoranthene	19.360	202	16631	0.384	ng/ul	98
20) Pyrene	19.719	202	16738	0.389	ng/ul	99
21) Benzo(a)anthracene	21.450	228	13194	0.410	ng/ul	99
22) Chrysene	21.503	228	14079	0.398	ng/ul	99
24) Benzo(b)fluoranthene	23.095	252	13831	0.417	ng/ul	98
25) Benzo(k)fluoranthene	23.143	252	13317	0.407	ng/ul	98
26) Benzo(a)pyrene	23.703	252	11452	0.416	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.201	276	13743	0.396	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.215	278	10953	0.395	ng/ul	95
29) Benzo(g,h,i)perylene	26.935	276	12090	0.388	ng/ul	93

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

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