

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032918.D
 Acq On : 09 Nov 2021 17:47
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4161

Quant Time: Nov 10 05:50:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 13:31:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.469	152	1731	0.400	ng/ul	0.00
4) Naphthalene-d8	10.244	136	6380	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.125	164	3877	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.865	188	8183	0.400	ng/ul	0.00
17) Chrysene-d12	21.057	240	6464	0.400	ng/ul	0.00
23) Perylene-d12	23.168	264	4949	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.867	96	896	0.405	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.845	152	3560	0.393	ng/ul	0.00
18) Fluoranthene-d10	18.896	212	8427	0.430	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.902	88	988	0.435	ng/ul	96
5) Naphthalene	10.293	128	7734	0.417	ng/ul	98
7) 2-Methylnaphthalene	11.917	142	5060	0.394	ng/ul	99
8) 1-Methylnaphthalene	12.142	142	5045	0.402	ng/ul	99
10) Acenaphthylene	13.842	152	6570	0.390	ng/ul	100
11) Acenaphthene	14.186	153	6009	0.408	ng/ul	99
12) Fluorene	15.179	166	7018	0.404	ng/ul	98
14) Pentachlorophenol	16.542	266	675	0.317	ng/ul	97
15) Phenanthrene	16.907	178	11055	0.407	ng/ul	100
16) Anthracene	16.997	178	9284	0.388	ng/ul	100
19) Fluoranthene	18.926	202	12712	0.419	ng/ul	99
20) Pyrene	19.289	202	12996	0.412	ng/ul	98
21) Benzo(a)anthracene	21.042	228	9307	0.390	ng/ul	99
22) Chrysene	21.093	228	11136	0.414	ng/ul	100
24) Benzo(b)fluoranthene	22.543	252	9600	0.423	ng/ul	98
25) Benzo(k)fluoranthene	22.584	252	10452	0.429	ng/ul	97
26) Benzo(a)pyrene	23.076	252	7620	0.407	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.232	276	9074	0.421	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.237	278	7226	0.424	ng/ul	99
29) Benzo(g,h,i)perylene	25.862	276	8111	0.434	ng/ul	100

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

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