

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032975.D
 Acq On : 11 Nov 2021 04:45
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4165

Quant Time: Nov 11 09:35:51 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	7.459	152	2300	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.235	136	8295	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.118	164	4692	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.859	188	9578	0.400	ng/ul	0.00
17) Chrysene-d12	21.047	240	7982	0.400	ng/ul	0.00
23) Perylene-d12	23.154	264	6489	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.864	96	1216	0.414	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.831	152	4504	0.382	ng/ul	0.00
18) Fluoranthene-d10	18.889	212	9949	0.411	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.898	88	1348	0.447	ng/ul#	93
5) Naphthalene	10.280	128	10131	0.421	ng/ul	98
7) 2-Methylnaphthalene	11.908	142	6449	0.387	ng/ul	99
8) 1-Methylnaphthalene	12.128	142	6373	0.390	ng/ul	100
10) Acenaphthylene	13.833	152	7931	0.389	ng/ul	100
11) Acenaphthene	14.178	153	7334	0.412	ng/ul	99
12) Fluorene	15.168	166	8190	0.390	ng/ul	99
14) Pentachlorophenol	16.536	266	928	0.372	ng/ul	99
15) Phenanthrene	16.900	178	12935	0.407	ng/ul	100
16) Anthracene	16.990	178	10730	0.384	ng/ul	100
19) Fluoranthene	18.919	202	15049	0.402	ng/ul	98
20) Pyrene	19.285	202	15449	0.397	ng/ul	100
21) Benzo(a)anthracene	21.033	228	11255	0.382	ng/ul	99
22) Chrysene	21.084	228	13568	0.408	ng/ul	100
24) Benzo(b)fluoranthene	22.531	252	12348	0.415	ng/ul	98
25) Benzo(k)fluoranthene	22.572	252	13316	0.417	ng/ul	99
26) Benzo(a)pyrene	23.064	252	9926	0.404	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.213	276	11759	0.416	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.218	278	9362	0.419	ng/ul	99
29) Benzo(g,h,i)perylene	25.843	276	10339	0.422	ng/ul	98

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

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