

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012422\
 Data File : BM033858.D
 Acq On : 25 Jan 2022 04:31
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4052

Quant Time: Jan 25 05:20:32 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	7.943	152	2143	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	7570	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.580	164	3721	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	7600	0.400	ng/ul	0.00
17) Chrysene-d12	21.480	240	6376	0.400	ng/ul #	0.00
23) Perylene-d12	23.878	264	6354	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	973	0.434	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	4214	0.404	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	8080	0.411	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.330	88	975	0.416	ng/ul#	81
5) Naphthalene	10.805	128	8365	0.402	ng/ul#	89
7) 2-Methylnaphthalene	12.416	142	5427	0.401	ng/ul	99
8) 1-Methylnaphthalene	12.636	142	5359	0.404	ng/ul	99
10) Acenaphthylene	14.303	152	6675	0.401	ng/ul#	91
11) Acenaphthene	14.640	153	5420	0.407	ng/ul	96
12) Fluorene	15.626	166	6034	0.397	ng/ul	97
14) Pentachlorophenol	16.972	266	1462	0.419	ng/ul	99
15) Phenanthrene	17.358	178	9762	0.403	ng/ul	95
16) Anthracene	17.451	178	8910	0.393	ng/ul	93
19) Fluoranthene	19.365	202	11842	0.407	ng/ul	99
20) Pyrene	19.727	202	12201	0.410	ng/ul	96
21) Benzo(a)anthracene	21.465	228	10294	0.395	ng/ul	97
22) Chrysene	21.514	228	11196	0.405	ng/ul	98
24) Benzo(b)fluoranthene	23.144	252	11382	0.399	ng/ul	91
25) Benzo(k)fluoranthene	23.192	252	11832	0.414	ng/ul	91
26) Benzo(a)pyrene	23.774	252	10342	0.418	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.383	276	13242	0.402	ng/ul#	84
28) Dibenzo(a,h)anthracene	26.407	278	10346	0.399	ng/ul	92
29) Benzo(g,h,i)perylene	27.151	276	11343	0.406	ng/ul	94

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

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