

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM036935.D
 Acq On : 11 Oct 2022 08:31
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4139

Quant Time: Oct 11 09:05:04 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM100622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 11 09:03:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	7689	0.400	ng/ul	0.00
4) Naphthalene-d8	10.754	136	26351	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.575	164	14678	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	29640	0.400	ng/ul	0.00
17) Chrysene-d12	21.485	240	20527	0.400	ng/ul	0.00
23) Perylene-d12	23.876	264	14512	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.348	96	3717	0.343	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.338	152	15774	0.396	ng/ul	0.00
18) Fluoranthene-d10	19.336	212	29292	0.477	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.386	88	3765	0.355	ng/ul#	85
5) Naphthalene	10.803	128	30228	0.401	ng/ul	96
7) 2-Methylnaphthalene	12.415	142	19062	0.414	ng/ul	91
8) 1-Methylnaphthalene	12.629	142	19329	0.413	ng/ul	99
10) Acenaphthylene	14.297	152	25237	0.414	ng/ul	98
11) Acenaphthene	14.640	153	21155	0.400	ng/ul	99
12) Fluorene	15.620	166	23368	0.390	ng/ul	99
14) Pentachlorophenol	16.967	266	2147	0.299	ng/ul	98
15) Phenanthrene	17.355	178	37619	0.398	ng/ul	99
16) Anthracene	17.448	178	32337	0.409	ng/ul	98
19) Fluoranthene	19.363	202	39711	0.486	ng/ul	100
20) Pyrene	19.726	202	40266	0.473	ng/ul	99
21) Benzo(a)anthracene	21.468	228	30728	0.414	ng/ul	98
22) Chrysene	21.520	228	31829	0.400	ng/ul	99
24) Benzo(b)fluoranthene	23.145	252	26950	0.416	ng/ul	95
25) Benzo(k)fluoranthene	23.195	252	25672	0.406	ng/ul	92
26) Benzo(a)pyrene	23.771	252	21480	0.408	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.363	276	27177	0.365	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.380	278	21607	0.359	ng/ul#	90
29) Benzo(g,h,i)perylene	27.128	276	23075	0.351	ng/ul	99

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

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