

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101022\
 Data File : BM036991.D
 Acq On : 12 Oct 2022 19:33
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 95 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4143

Quant Time: Oct 12 23:12:54 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	8594	0.400	ng/u1	0.00
4) Naphthalene-d8	10.754	136	31200	0.400	ng/u1	# 0.00
9) Acenaphthene-d10	14.575	164	16118	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.317	188	27403	0.400	ng/u1	0.00
17) Chrysene-d12	21.485	240	15329	0.400	ng/u1	0.00
23) Perylene-d12	23.882	264	14110	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.348	96	3891	0.321	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.338	152	18459	0.392	ng/u1	0.00
18) Fluoranthene-d10	19.336	212	23375	0.510	ng/u1	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.386	88	3948	0.333	ng/u1	90
5) Naphthalene	10.803	128	34865	0.391	ng/u1	97
7) 2-Methylnaphthalene	12.409	142	21646	0.397	ng/u1	92
8) 1-Methylnaphthalene	12.629	142	21903	0.395	ng/u1	99
10) Acenaphthylene	14.297	152	28564	0.426	ng/u1	98
11) Acenaphthene	14.640	153	22600	0.389	ng/u1	98
12) Fluorene	15.621	166	23781	0.361	ng/u1	98
14) Pentachlorophenol	16.971	266	1992	0.300	ng/u1	99
15) Phenanthrene	17.355	178	33511	0.383	ng/u1	99
16) Anthracene	17.448	178	30226	0.413	ng/u1	98
19) Fluoranthene	19.368	202	30398	0.499	ng/u1	98
20) Pyrene	19.726	202	30540	0.480	ng/u1	99
21) Benzo(a)anthracene	21.468	228	23107	0.417	ng/u1	100
22) Chrysene	21.520	228	22928	0.386	ng/u1	99
24) Benzo(b)fluoranthene	23.151	252	22264	0.353	ng/u1	96
25) Benzo(k)fluoranthene	23.198	252	22073	0.359	ng/u1	97
26) Benzo(a)pyrene	23.774	252	20025	0.392	ng/u1	94
27) Indeno(1,2,3-cd)pyrene	26.373	276	27036	0.373	ng/u1	99
28) Dibenzo(a,h)anthracene	26.390	278	21490	0.367	ng/u1	95
29) Benzo(g,h,i)perylene	27.138	276	22825	0.358	ng/u1	99

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

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