

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100722\
 Data File : BM036883.D
 Acq On : 08 Oct 2022 07:22
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4135

Quant Time: Oct 10 01:24:39 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 : Mon Oct 10 00:54:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.959	152	7644	0.400	ng/ul	0.00
4) Naphthalene-d8	10.765	136	26311	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.585	164	15064	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.323	188	31003	0.400	ng/ul	0.00
17) Chrysene-d12	21.490	240	24916	0.400	ng/ul	0.00
23) Perylene-d12	23.890	264	19002	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	3718	0.345	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.349	152	16109	0.405	ng/ul	0.00
18) Fluoranthene-d10	19.341	212	32858	0.441	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.390	88	3779	0.359	ng/ul#	87
5) Naphthalene	10.814	128	30200	0.401	ng/ul	96
7) 2-Methylnaphthalene	12.420	142	19260	0.419	ng/ul	91
8) 1-Methylnaphthalene	12.640	142	19658	0.421	ng/ul	98
10) Acenaphthylene	14.308	152	26350	0.421	ng/ul	98
11) Acenaphthene	14.645	153	21966	0.404	ng/ul	99
12) Fluorene	15.631	166	24710	0.401	ng/ul	99
14) Pentachlorophenol	16.976	266	3170	0.422	ng/ul	98
15) Phenanthrene	17.365	178	39990	0.404	ng/ul	99
16) Anthracene	17.453	178	34858	0.421	ng/ul	99
19) Fluoranthene	19.374	202	44434	0.448	ng/ul	97
20) Pyrene	19.732	202	45773	0.443	ng/ul	100
21) Benzo(a)anthracene	21.473	228	38561	0.428	ng/ul	99
22) Chrysene	21.525	228	39564	0.409	ng/ul	99
24) Benzo(b)fluoranthene	23.156	252	35228	0.415	ng/ul	95
25) Benzo(k)fluoranthene	23.203	252	34479	0.417	ng/ul	92
26) Benzo(a)pyrene	23.782	252	29071	0.422	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.382	276	33302	0.341	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.399	278	26332	0.334	ng/ul#	91
29) Benzo(g,h,i)perylene	27.147	276	27673	0.322	ng/ul	98

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

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