

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082123\
 Data File : BN026998.D
 Acq On : 21 Aug 2023 19:02
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4394

Quant Time: Aug 22 01:39:41 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 18 13:11:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.101	152	6596	0.400	ng/u1	0.00
4) Naphthalene-d8	10.927	136	20247	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.726	164	12399	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.472	188	27295	0.400	ng/u1	0.00
17) Chrysene-d12	21.647	240	21153	0.400	ng/u1	0.00
23) Perylene-d12	24.187	264	19860	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.439	96	3091	0.395	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.500	152	13035	0.438	ng/u1	0.00
18) Fluoranthene-d10	19.487	212	29851	0.447	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.477	88	3149	0.394	ng/u1	94
5) Naphthalene	10.977	128	22659	0.396	ng/u1	100
7) 2-Methylnaphthalene	12.577	142	15363	0.443	ng/u1	100
8) 1-Methylnaphthalene	12.791	142	16131	0.445	ng/u1	100
10) Acenaphthylene	14.458	152	23221	0.334	ng/u1	100
11) Acenaphthene	14.791	153	18484	0.384	ng/u1	99
12) Fluorene	15.771	166	20959	0.385	ng/u1	100
14) Pentachlorophenol	17.101	266	2227	0.288	ng/u1	99
15) Phenanthrene	17.515	178	34419	0.384	ng/u1	100
16) Anthracene	17.607	178	29198	0.360	ng/u1	100
19) Fluoranthene	19.520	202	39649	0.451	ng/u1	98
20) Pyrene	19.882	202	41077	0.438	ng/u1	100
21) Benzo(a)anthracene	21.627	228	32589	0.387	ng/u1	100
22) Chrysene	21.685	228	33906	0.408	ng/u1	99
24) Benzo(b)fluoranthene	23.395	252	33922	0.452	ng/u1	96
25) Benzo(k)fluoranthene	23.448	252	32546	0.439	ng/u1#	95
26) Benzo(a)pyrene	24.070	252	27000	0.398	ng/u1#	93
27) Indeno(1,2,3-cd)pyrene	26.871	276	31241	0.352	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.901	278	24240	0.343	ng/u1	97
29) Benzo(g,h,i)perylene	27.710	276	26690	0.362	ng/u1	98

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

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