

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020868.D
 Acq On : 26 Jul 2022 18:49
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4254

Quant Time: Jul 26 19:25:27 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 18:06:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	2439	0.400	ng/ul	0.00
4) Naphthalene-d8	10.612	136	8419	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.466	164	5183	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.218	188	10838	0.400	ng/ul	0.00
17) Chrysene-d12	21.421	240	9635	0.400	ng/ul	0.00
23) Perylene-d12	23.768	264	8092	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1097	0.374	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.212	152	5391	0.421	ng/ul	0.00
18) Fluoranthene-d10	19.254	212	12548	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	1146	0.381	ng/ul	96
5) Naphthalene	10.661	128	9177	0.348	ng/ul	99
7) 2-Methylnaphthalene	12.289	142	5926	0.355	ng/ul	100
8) 1-Methylnaphthalene	12.503	142	6571	0.407	ng/ul	100
10) Acenaphthylene	14.188	152	8991	0.349	ng/ul	100
11) Acenaphthene	14.530	153	7131	0.342	ng/ul	99
12) Fluorene	15.520	166	8038	0.351	ng/ul	99
14) Pentachlorophenol	16.884	266	728	0.446	ng/ul	98
15) Phenanthrene	17.260	178	13379	0.337	ng/ul	100
16) Anthracene	17.353	178	11101	0.349	ng/ul	99
19) Fluoranthene	19.286	202	15245	0.338	ng/ul	99
20) Pyrene	19.649	202	15732	0.334	ng/ul	100
21) Benzo(a)anthracene	21.406	228	11532	0.362	ng/ul	100
22) Chrysene	21.456	228	13664	0.333	ng/ul	99
24) Benzo(b)fluoranthene	23.055	252	12530	0.327	ng/ul	98
25) Benzo(k)fluoranthene	23.101	252	13257	0.334	ng/ul	98
26) Benzo(a)pyrene	23.666	252	11695	0.331	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.199	276	12274	0.330	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.219	278	9599	0.338	ng/ul	99
29) Benzo(g,h,i)perylene	26.937	276	10912	0.325	ng/ul	99

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

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