

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033023\
 Data File : BN024696.D
 Acq On : 01 Apr 2023 02:01
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4256

Quant Time: Apr 01 04:05:56 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 01 04:05:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.000	152	10887	0.400	ng/ul	0.00
4) Naphthalene-d8	10.812	136	33325	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.638	164	18653	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.384	188	38902	0.400	ng/ul	0.00
17) Chrysene-d12	21.565	240	32604	0.400	ng/ul	0.00
23) Perylene-d12	24.044	264	25946	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	4304	0.385	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.395	152	16291	0.396	ng/ul	0.00
18) Fluoranthene-d10	19.408	212	31424	0.394	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.405	88	4571	0.389	ng/ul	95
5) Naphthalene	10.861	128	32055	0.392	ng/ul	100
7) 2-Methylnaphthalene	12.467	142	20094	0.394	ng/ul	100
8) 1-Methylnaphthalene	12.687	142	21015	0.394	ng/ul	100
10) Acenaphthylene	14.356	152	25929	0.393	ng/ul	100
11) Acenaphthene	14.698	153	22049	0.385	ng/ul	100
12) Fluorene	15.683	166	24786	0.379	ng/ul	98
14) Pentachlorophenol	17.033	266	2683	0.319	ng/ul	99
15) Phenanthrene	17.422	178	40659	0.381	ng/ul	100
16) Anthracene	17.514	178	34130	0.383	ng/ul	100
19) Fluoranthene	19.436	202	44078	0.384	ng/ul	99
20) Pyrene	19.799	202	44779	0.383	ng/ul	99
21) Benzo(a)anthracene	21.548	228	39283	0.401	ng/ul	99
22) Chrysene	21.603	228	41139	0.382	ng/ul	99
24) Benzo(b)fluoranthene	23.284	252	40244	0.373	ng/ul	100
25) Benzo(k)fluoranthene	23.334	252	37869	0.365	ng/ul	99
26) Benzo(a)pyrene	23.939	252	33034	0.390	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.646	276	37398	0.370	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.666	278	29335	0.363	ng/ul	99
29) Benzo(g,h,i)perylene	27.444	276	30441	0.350	ng/ul	99

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

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