

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN083021\
 Data File : BN016114.D
 Acq On : 30 Aug 2021 16:43
 Operator : CG/JU
 Sample : SSTD0.410
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4210

Quant Time: Aug 30 17:16:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN083021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 30 17:16:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	1320	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	5479	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	3258	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	6762	0.400	ng/ul	0.00
17) Chrysene-d12	21.436	240	6885	0.400	ng/ul	0.00
23) Perylene-d12	23.827	264	6524	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	796	0.445	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	3406	0.413	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	7417	0.400	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.348	88	749	0.415	ng/ul	100
5) Naphthalene	10.710	128	6056	0.428	ng/ul	100
7) 2-Methylnaphthalene	12.322	142	4104	0.429	ng/ul	100
8) 1-Methylnaphthalene	12.542	142	3792	0.405	ng/ul	100
10) Acenaphthylene	14.219	152	5118	0.445	ng/ul	100
11) Acenaphthene	14.562	153	4151	0.418	ng/ul	100
12) Fluorene	15.547	166	4799	0.408	ng/ul	100
14) Pentachlorophenol	16.896	266	815	0.576	ng/ul	100
15) Phenanthrene	17.284	178	7806	0.418	ng/ul	100
16) Anthracene	17.377	178	6994	0.434	ng/ul	100
19) Fluoranthene	19.304	202	8936	0.415	ng/ul	100
20) Pyrene	19.666	202	9139	0.422	ng/ul	100
21) Benzo(a)anthracene	21.422	228	9169	0.428	ng/ul	100
22) Chrysene	21.474	228	9700	0.421	ng/ul	100
24) Benzo(b)fluoranthene	23.099	252	9791	0.402	ng/ul	100
25) Benzo(k)fluoranthene	23.146	252	9932	0.405	ng/ul	100
26) Benzo(a)pyrene	23.722	252	8681	0.424	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.297	276	11305	0.417	ng/ul	100
28) Dibenzo(a,h)anthracene	26.320	278	9442	0.414	ng/ul	100
29) Benzo(g,h,i)perylene	27.058	276	9600	0.411	ng/ul	100

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

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