

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN083021\
 Data File : BN016116.D
 Acq On : 30 Aug 2021 17:56
 Operator : CG/JU
 Sample : SSTD1.612
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6212

Quant Time: Aug 30 18:38:25 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	1577	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	6519	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	3818	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.247	188	7804	0.400	ng/ul	0.00
17) Chrysene-d12	21.432	240	8094	0.400	ng/ul	0.00
23) Perylene-d12	23.823	264	7576	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	3222	1.508	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	15185	1.547	ng/ul	0.00
18) Fluoranthene-d10	19.277	212	33023	1.515	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	3178	1.473	ng/ul#	84
5) Naphthalene	10.710	128	24998	1.484	ng/ul	97
7) 2-Methylnaphthalene	12.322	142	17238	1.515	ng/ul	100
8) 1-Methylnaphthalene	12.542	142	16934	1.519	ng/ul	100
10) Acenaphthylene	14.220	152	21641	1.605	ng/ul	97
11) Acenaphthene	14.563	153	17286	1.487	ng/ul	100
12) Fluorene	15.548	166	20246	1.470	ng/ul	98
14) Pentachlorophenol	16.897	266	3127	1.914	ng/ul	98
15) Phenanthrene	17.285	178	31705	1.472	ng/ul	99
16) Anthracene	17.378	178	29446	1.584	ng/ul	98
19) Fluoranthene	19.305	202	37809	1.495	ng/ul	98
20) Pyrene	19.667	202	37402	1.469	ng/ul	99
21) Benzo(a)anthracene	21.418	228	38599	1.532	ng/ul	99
22) Chrysene	21.470	228	39812	1.468	ng/ul	99
24) Benzo(b)fluoranthene	23.095	252	41482	1.467	ng/ul	94
25) Benzo(k)fluoranthene	23.142	252	41036	1.440	ng/ul	94
26) Benzo(a)pyrene	23.718	252	36112	1.519	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.296	276	48917	1.555	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.313	278	41203	1.556	ng/ul	97
29) Benzo(g,h,i)perylene	27.050	276	41075	1.515	ng/ul	97

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

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