

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007438.D
 Acq On : 27 Aug 2019 18:35
 Operator : HP/JU
 Sample : SSTD0.829
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.829

Quant Time: Aug 27 19:26:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 18:27:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	4159	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	17531	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	9028	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	21206	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	18629	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	19482	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	21039	0.68	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	51570	0.75	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	37928	0.752	ng/ul	99
5) 2-Methylnaphthalene	11.84	142	25808	0.681	ng/ul	99
7) Acenaphthylene	13.76	152	38670	0.739	ng/ul	99
8) Acenaphthene	14.11	153	27698	0.774	ng/ul	98
9) Fluorene	15.11	166	31206	0.719	ng/ul	98
11) Pentachlorophenol	16.47	266	3541	0.606	ng/ul	98
12) Phenanthrene	16.84	178	51476	0.762	ng/ul	99
13) Anthracene	16.93	178	49377	0.714	ng/ul	99
15) Fluoranthene	18.87	202	65500	0.780	ng/ul	100
17) Pyrene	19.24	202	67064	0.806	ng/ul	97
18) Benzo(a)anthracene	21.01	228	64341	0.786	ng/ul	99
19) Chrysene	21.06	228	66874	0.873	ng/ul	100
21) Benzo(b)fluoranthene	22.51	252	64722	0.819	ng/ul	97
22) Benzo(k)fluoranthene	22.55	252	68630	0.896	ng/ul#	95
23) Benzo(a)pyrene	23.04	252	64073	0.862	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.19	276	72949	0.910	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.21	278	57729	0.902	ng/ul	98
26) Benzo(g,h,i)perylene	25.82	276	61829	0.920	ng/ul	98

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

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