

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082318\
 Data File : BN002463.D
 Acq On : 23 Aug 2018 13:36
 Operator : JU/SJ
 Sample : SSTD3.213
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.213

Quant Time: Aug 23 14:06:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 13:07:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	1579	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	6783	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	4004	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	9088	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	10640	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	9275	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	31539	2.94	ng/ul	-0.02
14) Fluoranthene-d10	19.09	212	82212	2.99	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.49	128	49399	2.826	ng/ul 96
5) 2-Methylnaphthalene	12.12	142	35997	2.882	ng/ul 98
7) Acenaphthylene	14.02	152	54615	2.790	ng/ul 95
8) Acenaphthene	14.37	153	40244	2.796	ng/ul 96
9) Fluorene	15.36	166	47032	2.798	ng/ul# 92
11) Pentachlorophenol	16.72	266	4856	2.822	ng/ul 97
12) Phenanthrene	17.09	178	78660	2.901	ng/ul# 89
13) Anthracene	17.19	178	74612	2.903	ng/ul# 91
15) Fluoranthene	19.12	202	97622	3.032	ng/ul 97
17) Pyrene	19.49	202	93605	2.420	ng/ul# 94
18) Benzo(a)anthracene	21.24	228	95775	2.603	ng/ul# 85
19) Chrysene	21.29	228	95250	2.602	ng/ul 97
21) Benzo(b)fluoranthene	22.81	252	97281	2.891	ng/ul 82
22) Benzo(k)fluoranthene	22.85	252	98277	2.928	ng/ul# 84
23) Benzo(a)pyrene	23.38	252	91106	2.916	ng/ul# 94
24) Indeno(1,2,3-cd)pyrene	25.69	276	101703	2.898	ng/ul# 86
25) Dibenzo(a,h)anthracene	25.70	278	81097	2.898	ng/ul# 87
26) Benzo(g,h,i)perylene	26.37	276	83811	2.809	ng/ul 94

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

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