

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005155.D
 Acq On : 18 Apr 2019 17:06
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.459

Quant Time: Apr 19 05:25:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 11:57:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2970	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	10886	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	6147	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	14184	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	11341	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	10817	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	7148	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	15680	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.42	128	12148	0.404	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	8548	0.400	ng/ul	99
7) Acenaphthylene	13.95	152	11979	0.389	ng/ul	99
8) Acenaphthene	14.30	153	9221	0.400	ng/ul	100
9) Fluorene	15.29	166	10738	0.402	ng/ul	98
11) Pentachlorophenol	16.63	266	1365	0.397	ng/ul	99
12) Phenanthrene	17.03	178	17631	0.400	ng/ul	99
13) Anthracene	17.12	178	15838	0.393	ng/ul	100
15) Fluoranthene	19.06	202	19464	0.390	ng/ul	98
17) Pyrene	19.42	202	19147	0.392	ng/ul	99
18) Benzo(a)anthracene	21.19	228	16993	0.393	ng/ul	99
19) Chrysene	21.24	228	17996	0.405	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	16691	0.401	ng/ul	99
22) Benzo(k)fluoranthene	22.78	252	16294	0.398	ng/ul	99
23) Benzo(a)pyrene	23.29	252	14758	0.399	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.56	276	16554	0.404	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.58	278	13686	0.408	ng/ul	99
26) Benzo(g,h,i)perylene	26.21	276	14239	0.407	ng/ul	99

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

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