

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021119\
 Data File : BN004374.D
 Acq On : 11 Feb 2019 12:12
 Operator : JU/SJ
 Sample : SSTD1.606
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.606

Quant Time: Feb 11 12:44:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 11:59:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	1336	0.40	ng/ul	0.00
2) Naphthalene-d8	10.54	136	5301	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	3142	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	7979	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	10125	0.40	ng/ul	0.00
20) Perylene-d12	23.60	264	9602	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	13914	1.71	ng/ul	0.00
14) Fluoranthene-d10	19.17	212	41263	1.76	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.59	128	24430	1.636	ng/ul	98
5) 2-Methylnaphthalene	12.20	142	18099	1.746	ng/ul	100
7) Acenaphthylene	14.11	152	23798	1.512	ng/ul	99
8) Acenaphthene	14.45	153	20310	1.613	ng/ul	98
9) Fluorene	15.44	166	24440	1.709	ng/ul	99
11) Pentachlorophenol	16.78	266	3069	2.656	ng/ul	97
12) Phenanthrene	17.17	178	42452	1.637	ng/ul	99
13) Anthracene	17.27	178	37055	1.572	ng/ul	98
15) Fluoranthene	19.20	202	53438	1.711	ng/ul	99
17) Pyrene	19.56	202	52465	1.516	ng/ul	99
18) Benzo(a)anthracene	21.31	228	53403	1.552	ng/ul	100
19) Chrysene	21.37	228	58995	1.599	ng/ul	100
21) Benzo(b)fluoranthene	22.91	252	62447	1.714	ng/ul	96
22) Benzo(k)fluoranthene	22.96	252	63418	1.641	ng/ul	96
23) Benzo(a)pyrene	23.50	252	57138	1.649	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.89	276	76502	2.131	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	62882	2.135	ng/ul	97
26) Benzo(g,h,i)perylene	26.60	276	67292	2.228	ng/ul	99

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

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