

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053019\
 Data File : BN005958.D
 Acq On : 30 May 2019 14:24
 Operator : JU/SJ
 Sample : SSTD1.692
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.692

Quant Time: May 30 14:59:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 13:56:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	8227	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	30734	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	16612	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.15	188	37699	0.40	ng/ul	0.00
16) Chrysene-d12	21.36	240	25789	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	22674	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	77191	1.60	ng/ul	0.00
14) Fluoranthene-d10	19.18	212	168647	1.43	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.58	128	137455	1.690	ng/ul	99
5) 2-Methylnaphthalene	12.20	142	96770	1.742	ng/ul	100
7) Acenaphthylene	14.11	152	152618	1.825	ng/ul	99
8) Acenaphthene	14.46	153	104303	1.763	ng/ul	99
9) Fluorene	15.45	166	122154	1.767	ng/ul	99
11) Pentachlorophenol	16.80	266	17096	1.750	ng/ul	100
12) Phenanthrene	17.19	178	197026	1.918	ng/ul	99
13) Anthracene	17.28	178	193384	1.887	ng/ul	100
15) Fluoranthene	19.22	202	218811	1.539	ng/ul	99
17) Pyrene	19.58	202	219937	2.309	ng/ul	99
18) Benzo(a)anthracene	21.34	228	185059	2.081	ng/ul	98
19) Chrysene	21.39	228	172455	1.579	ng/ul	99
21) Benzo(b)fluoranthene	22.95	252	154420	2.265	ng/ul	90
22) Benzo(k)fluoranthene	23.00	252	149005	1.744	ng/ul#	92
23) Benzo(a)pyrene	23.53	252	142547	1.871	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.93	276	151771	1.781	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.95	278	121049	1.718	ng/ul	92
26) Benzo(g,h,i)perylene	26.63	276	127144	1.750	ng/ul	96

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

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