

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081219\
 Data File : BM022067.D
 Acq On : 13 Aug 2019 13:18
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.455

Quant Time: Aug 13 17:32:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 13 08:25:01 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	649	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	2869	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	1546	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	3352	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	3102	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	3443	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	1897	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	4136	0.37	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.46	128	3134	0.387	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	2127	0.394	ng/ul	100
7) Acenaphthylene	14.00	152	2672	0.371	ng/ul#	84
8) Acenaphthene	14.34	153	2381	0.404	ng/ul	98
9) Fluorene	15.34	166	2545	0.388	ng/ul	98
11) Pentachlorophenol	16.71	266	251	0.406	ng/ul#	83
12) Phenanthrene	17.08	178	3886	0.398	ng/ul	99
13) Anthracene	17.18	178	3431	0.412	ng/ul	98
15) Fluoranthene	19.10	202	4736	0.362	ng/ul	98
17) Pyrene	19.47	202	4955	0.389	ng/ul	95
18) Benzo(a)anthracene	21.22	228	4362	0.393	ng/ul	99
19) Chrysene	21.27	228	5439	0.381	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	4770	0.411	ng/ul	96
22) Benzo(k)fluoranthene	22.83	252	5316	0.405	ng/ul	96
23) Benzo(a)pyrene	23.35	252	4837	0.410	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	25.67	276	5754	0.408	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.67	278	4648	0.408	ng/ul	99
26) Benzo(g,h,i)perylene	26.34	276	5063	0.399	ng/ul	97

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

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