

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080719\
 Data File : BN007037.D
 Acq On : 08 Aug 2019 22:03
 Operator : HP/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.430

Quant Time: Aug 09 02:04:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	7397	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	28643	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	17689	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	44545	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	37048	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	33390	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	19322	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	53090	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	30857	0.374	ng/ul	98
5) 2-Methylnaphthalene	11.90	142	23070	0.373	ng/ul	100
7) Acenaphthylene	13.82	152	37150	0.362	ng/ul	99
8) Acenaphthene	14.17	153	26189	0.373	ng/ul	98
9) Fluorene	15.17	166	31315	0.368	ng/ul	100
11) Pentachlorophenol	16.52	266	3880	0.316	ng/ul#	81
12) Phenanthrene	16.90	178	53401	0.376	ng/ul	99
13) Anthracene	16.99	178	51812	0.357	ng/ul	99
15) Fluoranthene	18.93	202	64731	0.367	ng/ul	99
17) Pyrene	19.29	202	65450	0.396	ng/ul	98
18) Benzo(a)anthracene	21.06	228	59982	0.369	ng/ul	99
19) Chrysene	21.11	228	56631	0.372	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	49200	0.363	ng/ul	98
22) Benzo(k)fluoranthene	22.62	252	50012	0.381	ng/ul	98
23) Benzo(a)pyrene	23.11	252	45731	0.359	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.30	276	48638	0.354	ng/ul#	51
25) Dibenzo(a,h)anthracene	25.32	278	40084	0.365	ng/ul	97
26) Benzo(g,h,i)perylene	25.93	276	40281	0.350	ng/ul	97

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

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