

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071619\
 Data File : BN006855.D
 Acq On : 16 Jul 2019 08:21
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.429

Quant Time: Jul 16 08:58:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2312	0.40	ng/ul	0.01
2) Naphthalene-d8	10.37	136	8817	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4607	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9838	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	7579	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	8138	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4996	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	10306	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.42	128	9874	0.390	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	6103	0.357	ng/ul	100
7) Acenaphthylene	12.97	152	8481	0.357	ng/ul	99
8) Acenaphthene	14.31	153	7017	0.383	ng/ul	98
9) Fluorene	15.31	166	7287	0.363	ng/ul	98
11) Pentachlorophenol	16.72	266	330	0.225	ng/ul	95
12) Phenanthrene	17.05	178	11744	0.371	ng/ul	99
13) Anthracene	17.15	178	9102	0.333	ng/ul	99
15) Fluoranthene	19.09	202	13527	0.346	ng/ul	99
17) Pyrene	19.45	202	14221	0.411	ng/ul	99
18) Benzo(a)anthracene	21.22	228	10318	0.331	ng/ul	99
19) Chrysene	21.28	228	14413	0.415	ng/ul	99
21) Benzo(b)fluoranthene	22.81	252	12405	0.372	ng/ul	97
22) Benzo(k)fluoranthene	22.85	252	13931	0.399	ng/ul	98
23) Benzo(a)pyrene	23.38	252	12108	0.371	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.70	276	14383	0.435	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.70	278	11010	0.419	ng/ul	96
26) Benzo(g,h,i)perylene	26.38	276	13026	0.471	ng/ul	97

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

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