

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN051519\
 Data File : BN005695.D
 Acq On : 15 May 2019 09:02
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.477

Quant Time: May 15 11:59:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 14 07:15:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1179	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	5190	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	3547	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	9168	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	9880	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	10175	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3625	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	11695	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.40	128	5611	0.409	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	4226	0.451	ng/ul	100
7) Acenaphthylene	13.94	152	7360	0.412	ng/ul	100
8) Acenaphthene	14.28	153	5239	0.415	ng/ul	98
9) Fluorene	15.29	166	6240	0.423	ng/ul	99
11) Pentachlorophenol	16.68	266	726	0.306	ng/ul	90
12) Phenanthrene	17.02	178	10765	0.431	ng/ul	99
13) Anthracene	17.12	178	10338	0.415	ng/ul	99
15) Fluoranthene	19.06	202	14200	0.411	ng/ul	96
17) Pyrene	19.42	202	14806	0.406	ng/ul	96
18) Benzo(a)anthracene	21.20	228	14911	0.438	ng/ul	100
19) Chrysene	21.26	228	15928	0.381	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	14841	0.485	ng/ul	97
22) Benzo(k)fluoranthene	22.82	252	15654	0.408	ng/ul	99
23) Benzo(a)pyrene	23.34	252	14390	0.421	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.63	276	14398	0.376	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.64	278	11897	0.376	ng/ul	98
26) Benzo(g,h,i)perylene	26.30	276	11765	0.361	ng/ul	97

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

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