

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062719\
 Data File : BN006598.D
 Acq On : 27 Jun 2019 12:20
 Operator : HP/JU
 Sample : SSTD0.415
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Jun 27 14:09:12 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 : Thu Jun 27 14:04:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2576	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	10623	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	6014	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	13335	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11769	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	12108	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	6915	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	16277	0.45	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	12799	0.437	ng/ul	100
5) 2-Methylnaphthalene	12.08	142	8693	0.426	ng/ul	100
7) Acenaphthylene	14.00	152	13018	0.389	ng/ul	100
8) Acenaphthene	14.34	153	10134	0.437	ng/ul	100
9) Fluorene	15.34	166	11004	0.406	ng/ul	100
11) Pentachlorophenol	16.74	266	667	0.197	ng/ul	99
12) Phenanthrene	17.08	178	18074	0.422	ng/ul	100
13) Anthracene	17.17	178	15590	0.381	ng/ul	100
15) Fluoranthene	19.11	202	21990	0.465	ng/ul	100
17) Pyrene	19.48	202	22901	0.369	ng/ul	100
18) Benzo(a)anthracene	21.26	228	20334	0.391	ng/ul	100
19) Chrysene	21.32	228	22666	0.464	ng/ul	100
21) Benzo(b)fluoranthene	22.86	252	21242	0.421	ng/ul	100
22) Benzo(k)fluoranthene	22.90	252	21656	0.440	ng/ul	100
23) Benzo(a)pyrene	23.43	252	20241	0.430	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.77	276	20505	0.403	ng/ul	100
25) Dibenzo(a,h)anthracene	25.78	278	16232	0.399	ng/ul	100
26) Benzo(g,h,i)perylene	26.46	276	17209	0.404	ng/ul	100

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

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