

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006982.D
 Acq On : 07 Aug 2019 11:54
 Operator : HP/JU
 Sample : SSTD0.422
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.422

Quant Time: Aug 07 17:26:00 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 17:13:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	5237	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	20278	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	12800	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	32433	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	30019	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	32125	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	13925	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	40665	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.28	128	22699	0.389	ng/ul	100
5) 2-Methylnaphthalene	11.91	142	17084	0.390	ng/ul	100
7) Acenaphthylene	13.83	152	28752	0.388	ng/ul	100
8) Acenaphthene	14.18	153	19724	0.389	ng/ul	100
9) Fluorene	15.17	166	24007	0.390	ng/ul	100
11) Pentachlorophenol	16.52	266	3409	0.301	ng/ul	87
12) Phenanthrene	16.90	178	40250	0.389	ng/ul	100
13) Anthracene	16.99	178	41206	0.390	ng/ul	100
15) Fluoranthene	18.93	202	50344	0.392	ng/ul	100
17) Pyrene	19.30	202	51682	0.386	ng/ul	100
18) Benzo(a)anthracene	21.06	228	51164	0.388	ng/ul	100
19) Chrysene	21.11	228	48140	0.390	ng/ul	100
21) Benzo(b)fluoranthene	22.58	252	49600	0.381	ng/ul	100
22) Benzo(k)fluoranthene	22.62	252	49877	0.395	ng/ul	100
23) Benzo(a)pyrene	23.12	252	47451	0.387	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.31	276	51905	0.393	ng/ul	100
25) Dibenzo(a,h)anthracene	25.33	278	41527	0.393	ng/ul	100
26) Benzo(g,h,i)perylene	25.94	276	43605	0.393	ng/ul	100

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

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