

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101019\
 Data File : BN008063.D
 Acq On : 10 Oct 2019 15:16
 Operator : JU
 Sample : SSTD3.206
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.206

Quant Time: Oct 10 16:14:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 16:00:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2048	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	9954	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	6205	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	15455	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	16106	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	18405	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	53464	3.46	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	164235	3.27	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.45	128	84245	2.997	ng/ul 97
5) 2-Methylnaphthalene	12.07	142	61545	3.260	ng/ul 99
7) Acenaphthylene	13.98	152	96246	2.788	ng/ul 99
8) Acenaphthene	14.33	153	71333	2.875	ng/ul 99
9) Fluorene	15.32	166	85449	3.183	ng/ul 98
11) Pentachlorophenol	16.67	266	13106	3.764	ng/ul 99
12) Phenanthrene	17.05	178	141053	2.925	ng/ul 98
13) Anthracene	17.15	178	134954	2.969	ng/ul 98
15) Fluoranthene	19.08	202	183031	2.835	ng/ul 98
17) Pyrene	19.44	202	187546	2.380	ng/ul 98
18) Benzo(a)anthracene	21.20	228	189037	2.625	ng/ul 99
19) Chrysene	21.25	228	202239	2.616	ng/ul 98
21) Benzo(b)fluoranthene	22.75	252	199941	2.487	ng/ul 89
22) Benzo(k)fluoranthene	22.80	252	206379	2.455	ng/ul# 89
23) Benzo(a)pyrene	23.31	252	197246	2.462	ng/ul# 83
24) Indeno(1,2,3-cd)pyrene	25.59	276	236231	2.634	ng/ul 98
25) Dibenzo(a,h)anthracene	25.60	278	192155	2.707	ng/ul# 92
26) Benzo(g,h,i)perylene	26.24	276	208408	2.721	ng/ul 96

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

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