

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081220\
 Data File : BN011416.D
 Acq On : 12 Aug 2020 12:01
 Operator : CG/JU
 Sample : SST00.805
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.805

Quant Time: Aug 12 12:41:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 12 11:51:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	603	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	2133	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	1050	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	2562	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	2293	0.40	ng/ul	0.00
20) Perylene-d12	23.16	264	2243	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	2607	0.61	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	5194	0.46	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.24	128	4763	0.839	ng/ul	99
5) 2-Methylnaphthalene	11.87	142	3523	1.051	ng/ul	98
7) Acenaphthylene	13.79	152	4255	0.894	ng/ul	99
8) Acenaphthene	14.14	153	3361	0.889	ng/ul	99
9) Fluorene	15.14	166	3774	0.856	ng/ul	98
11) Pentachlorophenol	16.50	266	457	0.804	ng/ul	99
12) Phenanthrene	16.87	178	7128	0.906	ng/ul	99
13) Anthracene	16.97	178	5713	0.862	ng/ul	98
15) Fluoranthene	18.90	202	7054	0.725	ng/ul	99
17) Pyrene	19.27	202	7031	0.780	ng/ul	99
18) Benzo(a)anthracene	21.04	228	6498	0.907	ng/ul	99
19) Chrysene	21.09	228	8260	0.893	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	7240	0.855	ng/ul	98
22) Benzo(k)fluoranthene	22.58	252	7769	0.827	ng/ul	97
23) Benzo(a)pyrene	23.07	252	6373	0.840	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.22	276	8459	0.835	ng/ul	99
25) Dibenzo(a,h)anthracene	25.23	278	6937	0.841	ng/ul	98
26) Benzo(g,h,i)perylene	25.83	276	7538	0.772	ng/ul	99

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

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