

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111519\
 Data File : BN008591.D
 Acq On : 15 Nov 2019 18:06
 Operator : JU
 Sample : SSTD1.605
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6051

Quant Time: Nov 15 18:41:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 15 17:05:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2824	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	9538	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	5119	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	12466	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	13031	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	12277	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	25096	1.67	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	61600	1.66	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.43	128	45691	1.683	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	32307	1.787	ng/ul	99
7) Acenaphthylene	13.96	152	47104	1.975	ng/ul	98
8) Acenaphthene	14.31	153	34347	1.779	ng/ul	99
9) Fluorene	15.30	166	40609	1.921	ng/ul	99
11) Pentachlorophenol	16.65	266	7772	2.864	ng/ul	99
12) Phenanthrene	17.03	178	67826	1.790	ng/ul	98
13) Anthracene	17.12	178	64955	1.957	ng/ul	98
15) Fluoranthene	19.05	202	82941	1.909	ng/ul	98
17) Pyrene	19.41	202	82695	1.300	ng/ul	98
18) Benzo(a)anthracene	21.17	228	82358	1.815	ng/ul	99
19) Chrysene	21.22	228	79145	1.371	ng/ul	99
21) Benzo(b)fluoranthene	22.72	252	81057	1.772	ng/ul	92
22) Benzo(k)fluoranthene	22.76	252	78810	1.515	ng/ul#	89
23) Benzo(a)pyrene	23.27	252	76008	1.650	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.50	276	89239	1.704	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.52	278	72054	1.736	ng/ul#	93
26) Benzo(g,h,i)perylene	26.15	276	74695	1.592	ng/ul	96

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

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