

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111519\
 Data File : BN008593.D
 Acq On : 15 Nov 2019 19:17
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV071

Quant Time: Nov 15 22:44:14 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 22:41:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2291	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7504	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	4335	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	9643	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	10818	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	10094	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4953	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	11973	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	9047	0.400	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	6344	0.402	ng/ul	99
7) Acenaphthylene	13.96	152	8912	0.372	ng/ul	99
8) Acenaphthene	14.31	153	6762	0.380	ng/ul	100
9) Fluorene	15.30	166	7871	0.377	ng/ul	100
11) Pentachlorophenol	16.65	266	1360	0.373	ng/ul	99
12) Phenanthrene	17.03	178	13473	0.414	ng/ul	100
13) Anthracene	17.12	178	12409	0.404	ng/ul	100
15) Fluoranthene	19.05	202	16277	0.407	ng/ul	97
17) Pyrene	19.41	202	16631	0.383	ng/ul	98
18) Benzo(a)anthracene	21.17	228	16602	0.382	ng/ul	100
19) Chrysene	21.22	228	16724	0.391	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	16206	0.389	ng/ul	99
22) Benzo(k)fluoranthene	22.76	252	16485	0.399	ng/ul	99
23) Benzo(a)pyrene	23.26	252	15436	0.395	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.50	276	18048	0.395	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.52	278	14499	0.393	ng/ul	100
26) Benzo(g,h,i)perylene	26.15	276	15042	0.394	ng/ul	100

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

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