

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101019\
 Data File : BN008064.D
 Acq On : 10 Oct 2019 16:26
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV07

Quant Time: Oct 10 16:59:15 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 16:11:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1775	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.39	136	7947	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	5296	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	11384	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	12016	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	15176	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	5624	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	16865	0.43	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	9701	0.429	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	6507	0.419	ng/ul	100
7) Acenaphthylene	13.98	152	9969	0.395	ng/ul	100
8) Acenaphthene	14.33	153	7897	0.403	ng/ul	99
9) Fluorene	15.32	166	8880	0.399	ng/ul	98
11) Pentachlorophenol	16.68	266	1177	0.427	ng/ul	100
12) Phenanthrene	17.06	178	14467	0.434	ng/ul	99
13) Anthracene	17.15	178	12479	0.425	ng/ul	99
15) Fluoranthene	19.08	202	18978	0.431	ng/ul	100
17) Pyrene	19.44	202	20583	0.430	ng/ul	100
18) Benzo(a)anthracene	21.20	228	18215	0.417	ng/ul	99
19) Chrysene	21.26	228	23600	0.440	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	20313	0.411	ng/ul	100
22) Benzo(k)fluoranthene	22.80	252	23119	0.413	ng/ul	99
23) Benzo(a)pyrene	23.32	252	21353	0.411	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.59	276	25943	0.424	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	20511	0.424	ng/ul	99
26) Benzo(g,h,i)perylene	26.25	276	23908	0.426	ng/ul	99

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

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