

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050719\
 Data File : BN005495.D
 Acq On : 06 May 2019 21:03
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV64

Quant Time: May 07 04:11:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 03:05:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1661	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6120	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	3483	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8889	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	9854	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	11606	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3855	0.40	ng/ul	0.01
14) Fluoranthene-d10	19.04	212	11256	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	6446	0.398	ng/ul	100
5) 2-Methylnaphthalene	12.06	142	4413	0.399	ng/ul	98
7) Acenaphthylene	13.96	152	7048	0.402	ng/ul	100
8) Acenaphthene	14.31	153	5047	0.407	ng/ul	98
9) Fluorene	15.31	166	5879	0.406	ng/ul	98
11) Pentachlorophenol	16.68	266	808	0.351	ng/ul	89
12) Phenanthrene	17.04	178	9598	0.396	ng/ul	99
13) Anthracene	17.14	178	9607	0.398	ng/ul	99
15) Fluoranthene	19.07	202	13457	0.401	ng/ul	99
17) Pyrene	19.43	202	14260	0.392	ng/ul	100
18) Benzo(a)anthracene	21.21	228	13224	0.389	ng/ul	99
19) Chrysene	21.26	228	16705	0.400	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	13906	0.399	ng/ul	98
22) Benzo(k)fluoranthene	22.82	252	17628	0.403	ng/ul	99
23) Benzo(a)pyrene	23.33	252	15566	0.399	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.63	276	17363	0.398	ng/ul#	57
25) Dibenzo(a,h)anthracene	25.65	278	14349	0.398	ng/ul	100
26) Benzo(g,h,i)perylene	26.29	276	14633	0.393	ng/ul	99

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

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