

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053019\
 Data File : BN005960.D
 Acq On : 30 May 2019 15:43
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV94

Quant Time: May 30 16:20:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 15:28:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	7961	0.40	ng/ul	0.02
2) Naphthalene-d8	10.54	136	29173	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.41	164	15545	0.40	ng/ul	0.02
10) Phenanthrene-d10	17.16	188	34866	0.40	ng/ul	0.02
16) Chrysene-d12	21.38	240	21527	0.40	ng/ul	0.03
20) Perylene-d12	23.67	264	20693	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.14	152	18002	0.40	ng/ul	0.02
14) Fluoranthene-d10	19.20	212	35237	0.38	ng/ul	0.02
Target Compounds				Qvalue		
3) Naphthalene	10.59	128	32535	0.404	ng/ul	99
5) 2-Methylnaphthalene	12.21	142	22545	0.402	ng/ul	100
7) Acenaphthylene	14.13	152	34819	0.403	ng/ul	99
8) Acenaphthene	14.47	153	24296	0.405	ng/ul	99
9) Fluorene	15.46	166	27854	0.397	ng/ul	98
11) Pentachlorophenol	16.82	266	3266	0.369	ng/ul	99
12) Phenanthrene	17.20	178	44719	0.400	ng/ul	99
13) Anthracene	17.29	178	42948	0.402	ng/ul	100
15) Fluoranthene	19.23	202	46649	0.377	ng/ul	100
17) Pyrene	19.60	202	46595	0.411	ng/ul	99
18) Benzo(a)anthracene	21.36	228	37375	0.393	ng/ul	99
19) Chrysene	21.42	228	35485	0.397	ng/ul	99
21) Benzo(b)fluoranthene	22.98	252	34429	0.399	ng/ul	98
22) Benzo(k)fluoranthene	23.03	252	32829	0.391	ng/ul	100
23) Benzo(a)pyrene	23.57	252	32588	0.406	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.98	276	37028	0.426	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.99	278	29519	0.425	ng/ul	99
26) Benzo(g,h,i)perylene	26.69	276	31563	0.433	ng/ul	98

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

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