

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005150.D
 Acq On : 18 Apr 2019 12:00
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV58

Quant Time: Apr 18 12:34:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 11:57:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2826	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	10194	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	5493	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	12718	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	11261	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	11353	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	6445	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	14391	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.42	128	11191	0.397	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	7817	0.391	ng/ul	100
7) Acenaphthylene	13.95	152	10662	0.387	ng/ul	99
8) Acenaphthene	14.30	153	8191	0.397	ng/ul	99
9) Fluorene	15.30	166	9399	0.394	ng/ul	98
11) Pentachlorophenol	16.64	266	1114	0.361	ng/ul	98
12) Phenanthrene	17.03	178	15615	0.395	ng/ul	99
13) Anthracene	17.12	178	14091	0.390	ng/ul	100
15) Fluoranthene	19.06	202	17861	0.400	ng/ul	98
17) Pyrene	19.42	202	17709	0.365	ng/ul	100
18) Benzo(a)anthracene	21.18	228	16522	0.385	ng/ul	99
19) Chrysene	21.24	228	17474	0.396	ng/ul	98
21) Benzo(b)fluoranthene	22.74	252	16826	0.385	ng/ul	99
22) Benzo(k)fluoranthene	22.78	252	17105	0.398	ng/ul	99
23) Benzo(a)pyrene	23.29	252	15444	0.398	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.56	276	18061	0.420	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.58	278	14996	0.426	ng/ul	99
26) Benzo(g,h,i)perylene	26.21	276	15666	0.427	ng/ul	99

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

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