

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082318\
 Data File : BN002464.D
 Acq On : 23 Aug 2018 14:11
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV14

Quant Time: Aug 23 14:46:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 14:05:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	1235	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	5188	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	2969	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	7249	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	7896	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	7585	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	3285	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	8474	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.50	128	5265	0.394	ng/ul#	97
5) 2-Methylnaphthalene	12.13	142	3683	0.396	ng/ul	98
7) Acenaphthylene	14.03	152	5538	0.398	ng/ul	97
8) Acenaphthene	14.38	153	4185	0.395	ng/ul	94
9) Fluorene	15.37	166	4818	0.396	ng/ul#	92
11) Pentachlorophenol	16.74	266	311	0.318	ng/ul	90
12) Phenanthrene	17.10	178	8364	0.390	ng/ul#	89
13) Anthracene	17.20	178	7142	0.388	ng/ul#	92
15) Fluoranthene	19.13	202	9970	0.395	ng/ul	96
17) Pyrene	19.49	202	10446	0.388	ng/ul	96
18) Benzo(a)anthracene	21.25	228	9269	0.393	ng/ul#	85
19) Chrysene	21.30	228	10581	0.396	ng/ul	96
21) Benzo(b)fluoranthene	22.82	252	10744	0.394	ng/ul	87
22) Benzo(k)fluoranthene	22.86	252	11079	0.394	ng/ul#	87
23) Benzo(a)pyrene	23.39	252	10208	0.399	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.71	276	10939	0.410	ng/ul#	85
25) Dibenzo(a,h)anthracene	25.72	278	8609	0.409	ng/ul	91
26) Benzo(g,h,i)perylene	26.39	276	9344	0.409	ng/ul	95

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

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