

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN082819\
 Data File : BN007441.D
 Acq On : 27 Aug 2019 20:41
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV32

Quant Time: Aug 27 21:12:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 20:30:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3695	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	15477	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	7811	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	17811	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	15147	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	16143	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	9258	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	21988	0.38	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.21	128	16996	0.389	ng/ul	99
5) 2-Methylnaphthalene	11.84	142	11394	0.388	ng/ul	99
7) Acenaphthylene	13.76	152	16763	0.386	ng/ul	100
8) Acenaphthene	14.11	153	12136	0.389	ng/ul	99
9) Fluorene	15.11	166	13014	0.385	ng/ul	99
11) Pentachlorophenol	16.48	266	1345	0.335	ng/ul	98
12) Phenanthrene	16.84	178	21488	0.387	ng/ul	99
13) Anthracene	16.94	178	19986	0.382	ng/ul	99
15) Fluoranthene	18.87	202	28053	0.377	ng/ul	99
17) Pyrene	19.24	202	29226	0.394	ng/ul	97
18) Benzo(a)anthracene	21.01	228	25870	0.382	ng/ul	99
19) Chrysene	21.06	228	28706	0.395	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	26862	0.381	ng/ul	99
22) Benzo(k)fluoranthene	22.55	252	29414	0.399	ng/ul#	96
23) Benzo(a)pyrene	23.04	252	27163	0.386	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.19	276	30640	0.389	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.21	278	24136	0.388	ng/ul	99
26) Benzo(g,h,i)perylene	25.82	276	25992	0.387	ng/ul	100

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

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