

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004054.D
 Acq On : 18 Dec 2018 02:03
 Operator : JU/SJ
 Sample : J6270-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E01

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:41 PM

Quant Time: Dec 18 07:05:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 18 06:35:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1578	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7584	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4670	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12429m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14701m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13566m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3464	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	12512m	0.34	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	8023	0.382	ng/ul	98
8) Acenaphthene	14.53	153	565	0.031	ng/ul#	93
9) Fluorene	15.52	166	1031	0.048	ng/ul#	96
12) Phenanthrene	17.26	178	7014m	0.172	ng/ul	
13) Anthracene	17.35	178	8824m	0.263	ng/ul	
15) Fluoranthene	19.27	202	38470m	0.799	ng/ul	
17) Pyrene	19.64	202	39722m	0.753	ng/ul	
18) Benzo(a)anthracene	21.38	228	27408	0.590	ng/ul	92
19) Chrysene	21.43	228	32918m	0.619	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	54529m	0.972	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	18335m	0.335	ng/ul	
23) Benzo(a)pyrene	23.62	252	26555m	0.546	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	25598m	0.459	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	8679m	0.194	ng/ul	
26) Benzo(g,h,i)perylene	26.82	276	35068m	0.724	ng/ul	

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

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