

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002488.D
 Acq On : 24 Aug 2018 04:22
 Operator : JU/SJ
 Sample : J4424-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YS7

Manual Integrations
 APPROVED

Sohil
 8/24/2018 5:11:59 PM

Quant Time: Aug 24 06:54:16 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 : Fri Aug 24 02:20:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1869	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	7802	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.31	164	4433	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	10516m	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	11512m	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	10484m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	2362	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6746m	0.22	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.49	128	998	0.050	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	440	0.031	ng/ul	100
7) Acenaphthylene	14.02	152	7510	0.362	ng/ul	97
8) Acenaphthene	14.37	153	675	0.043	ng/ul#	93
9) Fluorene	15.36	166	1232	0.068	ng/ul	92
12) Phenanthrene	17.09	178	23496m	0.755	ng/ul	
13) Anthracene	17.18	178	6168m	0.231	ng/ul	
15) Fluoranthene	19.12	202	60899m	1.665	ng/ul	
17) Pyrene	19.48	202	68431	1.743	ng/ul#	95
18) Benzo(a)anthracene	21.23	228	31073	0.903	ng/ul#	84
19) Chrysene	21.28	228	38661	0.993	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	57349m	1.523	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	18523m	0.476	ng/ul	
23) Benzo(a)pyrene	23.37	252	37376	1.056	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.68	276	26057	0.706	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.68	278	6407	0.220	ng/ul	97
26) Benzo(a,h,i)perylene	26.36	276	24993	0.791	ng/ul	95

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

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