

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004065.D
 Acq On : 18 Dec 2018 08:29
 Operator : JU/SJ
 Sample : J6270-13
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E12

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:57:19 PM

Quant Time: Dec 18 11:18:45 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	7.83	152	1988	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9896	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5977	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14948m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16271m	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	16957	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4124	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	13020m	0.30	ng/ul	0.00

Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	13235	0.492	ng/ul	99
12) Phenanthrene	17.26	178	5224m	0.106	ng/ul	
13) Anthracene	17.35	178	12065m	0.299	ng/ul	
15) Fluoranthene	19.28	202	60383m	1.042	ng/ul	
17) Pyrene	19.64	202	65594	1.124	ng/ul	99
18) Benzo(a)anthracene	21.39	228	47838	0.931	ng/ul	97
19) Chrysene	21.44	228	50607	0.860	ng/ul	97
21) Benzo(b)fluoranthene	23.02	252	93832m	1.338	ng/ul	
22) Benzo(k)fluoranthene	23.07	252	30500m	0.446	ng/ul	
23) Benzo(a)pyrene	23.63	252	44016	0.725	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.09	276	39433	0.566	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.09	278	10974	0.196	ng/ul#	69
26) Benzo(g,h,i)perylene	26.82	276	36408	0.601	ng/ul	94

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

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