

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004072.D
 Acq On : 18 Dec 2018 12:36
 Operator : JU/SJ
 Sample : J6236-12DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E22DL

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:15:55 PM

Quant Time: Dec 19 02:53:58 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 : Wed Dec 19 01:54:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1641	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8145	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4931	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12390	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13572	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12743	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	506	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	1594	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	13634	0.614	ng/ul	99
12) Phenanthrene	17.26	178	3147	0.077	ng/ul	95
13) Anthracene	17.35	178	12789	0.383	ng/ul	100
15) Fluoranthene	19.28	202	57585	1.199	ng/ul	98
17) Pyrene	19.64	202	66792	1.372	ng/ul	96
18) Benzo(a)anthracene	21.39	228	44122	1.029	ng/ul	97
19) Chrysene	21.44	228	44046	0.898	ng/ul	97
21) Benzo(b)fluoranthene	23.02	252	79295m	1.504	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	27218m	0.530	ng/ul	
23) Benzo(a)pyrene	23.62	252	35243	0.772	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.09	276	34294	0.655	ng/ul	94
25) Dibenzo(a,h)anthracene	26.09	278	8141m	0.193	ng/ul	
26) Benzo(g,h,i)perylene	26.82	276	28132	0.618	ng/ul	97

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

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