

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121018\
 Data File : BN003790.D
 Acq On : 10 Dec 2018 22:24
 Operator : JU/SJ
 Sample : J6077-11DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y00DL

Manual Integrations
 APPROVED

Sohil
 12/11/2018 6:12:20 PM

Quant Time: Dec 11 04:40:44 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 11 04:22:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	5238	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	21200	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	10784	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	23515	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	20883	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	19494	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2251	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	4909	0.07	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.25	178	21836	0.283	ng/ul	99
13) Anthracene	17.34	178	1515	0.024	ng/ul#	87
15) Fluoranthene	19.27	202	74494	0.817	ng/ul	98
17) Pyrene	19.63	202	64623	0.863	ng/ul	97
18) Benzo(a)anthracene	21.38	228	21432	0.325	ng/ul	96
19) Chrysene	21.43	228	33999	0.450	ng/ul	98
21) Benzo(b)fluoranthene	23.00	252	45252m	0.561	ng/ul	
22) Benzo(k)fluoranthene	23.05	252	16145m	0.206	ng/ul	
23) Benzo(a)pyrene	23.60	252	21418	0.307	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.04	276	19051	0.238	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.05	278	3902	0.061	ng/ul#	70
26) Benzo(g,h,i)perylene	26.77	276	16126	0.232	ng/ul	96

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

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