

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121018\
 Data File : BN003792.D
 Acq On : 10 Dec 2018 23:34
 Operator : JU/SJ
 Sample : J6077-12DL 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y01DL

Manual Integrations
 APPROVED

Sohil
 12/11/2018 6:12:23 PM

Quant Time: Dec 11 04:47:37 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	5621	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	22889	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	11620	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	25218	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	22968	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	21334	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2011	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	4467	0.06	ng/ul	0.00
Target Compounds						Ovalue
12) Phenanthrene	17.25	178	14601	0.176	ng/ul	99
15) Fluoranthene	19.27	202	46465	0.475	ng/ul	98
17) Pyrene	19.63	202	40854	0.496	ng/ul	98
18) Benzo(a)anthracene	21.38	228	14044	0.194	ng/ul	96
19) Chrysene	21.43	228	20539	0.247	ng/ul	96
21) Benzo(b)fluoranthene	23.00	252	26898m	0.305	ng/ul	
22) Benzo(k)fluoranthene	23.04	252	9019m	0.105	ng/ul	
23) Benzo(a)pyrene	23.60	252	12494	0.163	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.04	276	10464	0.119	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.05	278	2235	0.032	ng/ul#	42
26) Benzo(g,h,i)perylene	26.77	276	9293	0.122	ng/ul#	92

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

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