

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003386.D
 Acq On : 01 Nov 2018 19:35
 Operator : MJ/SJ
 Sample : J5657-18MSD
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40F1MSD

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:07 PM

Quant Time: Nov 12 14:38:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 02 04:49:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	11596	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	47799	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	26944	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	65638	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	61711	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	51354	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	20343	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	55657	0.30	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.55	153	28791	0.275	ng/ul	99
11) Pentachlorophenol	16.87	266	136	0.010	ng/ul	100
12) Phenanthrene	17.27	178	20415	0.095	ng/ul	98
15) Fluoranthene	19.29	202	54038	0.223	ng/ul	94
17) Pyrene	19.65	202	130661	0.582	ng/ul	96
18) Benzo(a)anthracene	21.39	228	20148	0.104	ng/ul#	90
19) Chrysene	21.45	228	24989	0.114	ng/ul#	94
21) Benzo(b)fluoranthene	23.02	252	29964m	0.135	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	8422m	0.039	ng/ul	
23) Benzo(a)pyrene	23.62	252	18451	0.100	ng/ul#	62
24) Indeno(1,2,3-cd)pyrene	26.08	276	11037	0.056	ng/ul#	59
26) Benzo(g,h,i)perylene	26.81	276	9944	0.059	ng/ul#	76

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

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