

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE083118\
 Data File : BE097402.D
 Acq On : 31 Aug 2018 11:55
 Operator : SJ/JU
 Sample : J4596-06
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z29

Manual Integrations
 APPROVED

Sohil
 9/4/2018 4:38:31 PM

Quant Time: Aug 31 12:51:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 11:16:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	806	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4058	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2726	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	8510m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11411m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	10218m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1280	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	5313m	0.19	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.17	128	205	0.021	ng/ul#	52
12) Phenanthrene	16.80	178	1033m	0.047	ng/ul	
15) Fluoranthene	18.83	202	1947m	0.069	ng/ul	
17) Pyrene	19.20	202	2329m	0.076	ng/ul	
18) Benzo(a)anthracene	20.95	228	1144m	0.039	ng/ul	
19) Chrysene	20.98	228	2429m	0.072	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	2502m	0.091	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	1031m	0.034	ng/ul	
23) Benzo(a)pyrene	23.10	252	790	0.031	ng/ul#	50
24) Indeno(1,2,3-cd)pyrene	25.49	276	882	0.027	ng/ul#	47
26) Benzo(g,h,i)perylene	26.19	276	963	0.034	ng/ul#	72

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

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