

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097238.D
 Acq On : 8 Aug 2018 16:55
 Operator : SJ/JU
 Sample : J4268-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA1

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:43:16 PM

Quant Time: Aug 09 01:12:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 23:25:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1235	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6613	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3990	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9269m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	9511m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8979m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2005	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	5575m	0.17	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	866	0.054	ng/ul#	85
8) Acenaphthene	14.08	153	259	0.021	ng/ul#	88
9) Fluorene	15.08	166	320	0.022	ng/ul#	90
12) Phenanthrene	16.81	178	8139m	0.344	ng/ul	
13) Anthracene	16.90	178	1984m	0.090	ng/ul	
15) Fluoranthene	18.84	202	27057m	0.866	ng/ul	
17) Pyrene	19.21	202	24953	0.979	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	13823	0.509	ng/ul#	86
19) Chrysene	21.01	228	16147	0.577	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	22018m	0.900	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	6287m	0.238	ng/ul	
23) Benzo(a)pyrene	23.11	252	15865	0.648	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.50	276	10362	0.338	ng/ul#	72
25) Dibenzo(a,h)anthracene	25.52	278	2503	0.100	ng/ul#	92
26) Benzo(g,h,i)perylene	26.21	276	9755	0.368	ng/ul#	93

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

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