

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080918\
 Data File : BE097247.D
 Acq On : 9 Aug 2018 12:42
 Operator : SJ/JU
 Sample : J4268-13DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC3DL

Manual Integrations
 APPROVED

Sohil
 8/10/2018 2:50:21 PM

Quant Time: Aug 09 16:21:28 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	1271	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6782	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3863	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9139	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9850	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	10430	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	438	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	1076	0.03	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	753	0.046	ng/ul#	86
5) 2-Methylnaphthalene	11.80	142	382	0.035	ng/ul	98
7) Acenaphthylene	13.73	152	1507	0.090	ng/ul	93
8) Acenaphthene	14.08	153	669	0.055	ng/ul#	90
9) Fluorene	15.08	166	940	0.068	ng/ul	93
12) Phenanthrene	16.81	178	30742	1.320	ng/ul#	89
13) Anthracene	16.90	178	4393	0.202	ng/ul#	94
15) Fluoranthene	18.84	202	103060	3.347	ng/ul	84
17) Pyrene	19.20	202	91634	3.472	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	46127	1.639	ng/ul#	85
19) Chrysene	21.01	228	59408	2.048	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	85331m	3.002	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	26335m	0.859	ng/ul	
23) Benzo(a)pyrene	23.11	252	52670	1.852	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.50	276	43073	1.210	ng/ul#	71
25) Dibenzo(a,h)anthracene	25.52	278	9621	0.330	ng/ul	90
26) Benzo(a,h,i)perylene	26.20	276	41569	1.351	ng/ul#	91

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

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