

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101218\
 Data File : BE097897.D
 Acq On : 13 Oct 2018 2:11
 Operator : SJ/JU
 Sample : J5183-21DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC38DL

Manual Integrations
 APPROVED

Sohil
 10/15/2018 3:56:13 PM

Quant Time: Oct 13 04:27:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 13 02:57:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	2161	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11209	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	7300	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.32	188	17090	0.40	ng/ul	0.00
16) Chrysene-d12	21.50	240	17746	0.40	ng/ul	0.00
20) Perylene-d12	24.07	264	18368	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	453	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.34	212	1338	0.02	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	2643	0.097	ng/ul#	91
5) 2-Methylnaphthalene	12.37	142	3375	0.171	ng/ul	100
8) Acenaphthene	14.62	153	1359	0.063	ng/ul	97
9) Fluorene	15.61	166	1781	0.064	ng/ul#	92
12) Phenanthrene	17.36	178	28468	0.704	ng/ul	97
13) Anthracene	17.45	178	6688	0.178	ng/ul#	93
15) Fluoranthene	19.37	202	48250	0.929	ng/ul	95
17) Pyrene	19.74	202	55004	1.040	ng/ul	98
18) Benzo(a)anthracene	21.48	228	31496	0.587	ng/ul#	87
19) Chrysene	21.54	228	23918	0.496	ng/ul#	87
21) Benzo(b)fluoranthene	23.29	252	27948m	0.569	ng/ul	
22) Benzo(k)fluoranthene	23.33	252	7526m	0.143	ng/ul	
23) Benzo(a)pyrene	23.95	252	22214	0.461	ng/ul#	56
24) Indeno(1,2,3-cd)pyrene	26.79	276	12456	0.228	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.80	278	3647	0.081	ng/ul#	1
26) Benzo(g,h,i)perylene	27.63	276	18343	0.385	ng/ul#	74

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

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