

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098101.D
 Acq On : 26 Oct 2018 22:47
 Operator : SJ/JU
 Sample : J5598-12DL 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD6DL

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:32 PM

Quant Time: Oct 27 00:55:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 27 00:33:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1087	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5561	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4004	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	9602	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	13130	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	11125	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	261	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	769	0.02	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.72	128	286	0.021	ng/ul#	54
7) Acenaphthylene	14.25	152	771	0.049	ng/ul#	87
12) Phenanthrene	17.33	178	922	0.042	ng/ul#	90
13) Anthracene	17.42	178	543	0.025	ng/ul#	74
15) Fluoranthene	19.35	202	19031	0.674	ng/ul	99
17) Pyrene	19.71	202	20245	0.660	ng/ul	99
18) Benzo(a)anthracene	21.45	228	16046	0.416	ng/ul	98
19) Chrysene	21.51	228	16460	0.551	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	24144m	0.783	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	7549m	0.250	ng/ul	
23) Benzo(a)pyrene	23.91	252	10737	0.362	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.71	276	7203	0.212	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.73	278	2070	0.073	ng/ul#	69
26) Benzo(g,h,i)perylene	27.54	276	5625	0.197	ng/ul#	97

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

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