

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098094.D
 Acq On : 26 Oct 2018 18:30
 Operator : SJ/JU
 Sample : J5598-04DL 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE7DL

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:25 PM

Quant Time: Oct 27 00:39:45 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	965	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4776	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3354	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8029	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11504	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9465	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	415	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	1240	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	5335	0.467	ng/ul	99
5) 2-Methylnaphthalene	12.34	142	3221	0.373	ng/ul	100
7) Acenaphthylene	14.26	152	9357	0.707	ng/ul	99
8) Acenaphthene	14.60	153	504	0.051	ng/ul#	83
9) Fluorene	15.59	166	704	0.059	ng/ul#	94
12) Phenanthrene	17.33	178	42837	2.312	ng/ul	99
13) Anthracene	17.42	178	9941	0.543	ng/ul	99
15) Fluoranthene	19.35	202	323360	13.687	ng/ul	99
17) Pyrene	19.71	202	292094	10.874	ng/ul	98
18) Benzo(a)anthracene	21.45	228	186474	5.513	ng/ul	99
19) Chrysene	21.51	228	231331	8.841	ng/ul	99
21) Benzo(b)fluoranthene	23.25	252	321022m	12.242	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	76640m	2.986	ng/ul	
23) Benzo(a)pyrene	23.91	252	109586	4.345	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.71	276	90844	3.143	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.73	278	24472	1.016	ng/ul	98
26) Benzo(g,h,i)perylene	27.54	276	76138	3.129	ng/ul	96

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

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