

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098089.D
 Acq On : 26 Oct 2018 14:54
 Operator : SJ/JU
 Sample : J5598-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD7

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:22 PM

Quant Time: Oct 27 00:21:41 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 : Fri Oct 26 23:15:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1049	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5342	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3805	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	9046m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	10017m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	10319m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1700	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	5599m	0.19	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.72	128	1473	0.115	ng/ul#	90
5) 2-Methylnaphthalene	12.34	142	422	0.044	ng/ul	98
7) Acenaphthylene	14.26	152	11809	0.787	ng/ul	97
8) Acenaphthene	14.60	153	974	0.087	ng/ul	93
9) Fluorene	15.59	166	1403	0.104	ng/ul#	98
12) Phenanthrene	17.33	178	49089m	2.351	ng/ul	
13) Anthracene	17.42	178	10332m	0.501	ng/ul	
15) Fluoranthene	19.35	202	342803m	12.878	ng/ul	
17) Pyrene	19.71	202	307011	13.126	ng/ul	98
18) Benzo(a)anthracene	21.45	228	232027m	7.878	ng/ul	
19) Chrysene	21.51	228	215320	9.451	ng/ul	99
21) Benzo(b)fluoranthene	23.25	252	318956m	11.156	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	87975m	3.144	ng/ul	
23) Benzo(a)pyrene	23.91	252	118778m	4.320	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.72	276	95887m	3.043	ng/ul	
25) Dibenzo(a,h)anthracene	26.73	278	25270m	0.962	ng/ul	
26) Benzo(g,h,i)perylene	27.55	276	69905m	2.635	ng/ul	

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

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