

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098078.D
 Acq On : 26 Oct 2018 5:05
 Operator : SJ/JU
 Sample : J5598-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD5

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:47:56 PM

Quant Time: Oct 26 08:24:35 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 07:43:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1182	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5903	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	4383	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	10312m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11483m	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	11063m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2554	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	9172m	0.27	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	5052	0.358	ng/ul	100
5) 2-Methylnaphthalene	12.35	142	1414	0.132	ng/ul	100
7) Acenaphthylene	14.26	152	3802	0.220	ng/ul	96
8) Acenaphthene	14.60	153	679	0.053	ng/ul#	90
9) Fluorene	15.59	166	549	0.035	ng/ul#	93
12) Phenanthrene	17.33	178	9245m	0.388	ng/ul	
13) Anthracene	17.42	178	6148m	0.262	ng/ul	
15) Fluoranthene	19.35	202	72947m	2.404	ng/ul	
17) Pyrene	19.71	202	111704	4.166	ng/ul	99
18) Benzo(a)anthracene	21.46	228	82741m	2.451	ng/ul	
19) Chrysene	21.51	228	76944m	2.946	ng/ul	
21) Benzo(b)fluoranthene	23.24	252	176343m	5.753	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	49980m	1.666	ng/ul	
23) Benzo(a)pyrene	23.91	252	66180m	2.245	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.71	276	54300m	1.607	ng/ul	
25) Dibenzo(a,h)anthracene	26.73	278	13469m	0.478	ng/ul	
26) Benzo(g,h,i)perylene	27.55	276	39294m	1.382	ng/ul	

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

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