

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098076.D
 Acq On : 26 Oct 2018 3:51
 Operator : SJ/JU
 Sample : J5598-05
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE8

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:47:53 PM

Quant Time: Oct 26 08:13:14 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	7.88	152	1385	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	6854	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	4957	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	11688m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	12906m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	11871m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2324	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	7210m	0.19	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	8578	0.523	ng/ul	98
5) 2-Methylnaphthalene	12.35	142	1673	0.135	ng/ul	100
7) Acenaphthylene	14.25	152	7366	0.377	ng/ul	99
8) Acenaphthene	14.60	153	2079	0.142	ng/ul	97
9) Fluorene	15.59	166	1007	0.057	ng/ul#	95
12) Phenanthrene	17.33	178	20625m	0.765	ng/ul	
13) Anthracene	17.42	178	10368m	0.389	ng/ul	
15) Fluoranthene	19.35	202	171470m	4.986	ng/ul	
17) Pyrene	19.72	202	209841	6.964	ng/ul	97
18) Benzo(a)anthracene	21.45	228	191715m	5.052	ng/ul	
19) Chrysene	21.51	228	135477m	4.615	ng/ul	
21) Benzo(b)fluoranthene	23.25	252	292487m	8.893	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	91690m	2.848	ng/ul	
23) Benzo(a)pyrene	23.91	252	119561m	3.780	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.72	276	98517m	2.718	ng/ul	
25) Dibenzo(a,h)anthracene	26.73	278	27190m	0.900	ng/ul	
26) Benzo(g,h,i)perylene	27.55	276	67722m	2.219	ng/ul	

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

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