

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098087.D
 Acq On : 26 Oct 2018 13:39
 Operator : SJ/JU
 Sample : J5598-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE6

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:20 PM

Quant Time: Oct 27 00:07:09 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	1076	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5287	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3731	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8817m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	9898m	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9871m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1745	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	5461m	0.19	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	19579	1.547	ng/ul	95
5) 2-Methylnaphthalene	12.34	142	6448	0.674	ng/ul	100
7) Acenaphthylene	14.26	152	16655	1.131	ng/ul	99
8) Acenaphthene	14.60	153	1587	0.144	ng/ul	94
9) Fluorene	15.59	166	1304	0.099	ng/ul#	92
12) Phenanthrene	17.33	178	42817m	2.104	ng/ul	
13) Anthracene	17.42	178	31952m	1.590	ng/ul	
15) Fluoranthene	19.35	202	219252m	8.451	ng/ul	
17) Pyrene	19.71	202	276377	11.959	ng/ul	98
18) Benzo(a)anthracene	21.46	228	260783m	8.961	ng/ul	
19) Chrysene	21.51	228	290967	12.925	ng/ul	99
21) Benzo(b)fluoranthene	23.25	252	560804m	20.506	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	184461m	6.890	ng/ul	
23) Benzo(a)pyrene	23.91	252	167435m	6.366	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.72	276	218612m	7.253	ng/ul	
25) Dibenzo(a,h)anthracene	26.74	278	54312m	2.162	ng/ul	
26) Benzo(g,h,i)perylene	27.55	276	184038m	7.253	ng/ul	

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

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