

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098092.D
 Acq On : 26 Oct 2018 17:14
 Operator : SJ/JU
 Sample : J5598-11DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD5DL

Manual Integrations
 APPROVED

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

Quant Time: Oct 27 00:29:49 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	1023	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5149	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3650	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8770	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11685	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	10544	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	471	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	1637	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	974	0.079	ng/ul#	85
5) 2-Methylnaphthalene	12.34	142	291	0.031	ng/ul	96
7) Acenaphthylene	14.25	152	671	0.047	ng/ul#	78
12) Phenanthrene	17.33	178	1724	0.085	ng/ul#	93
13) Anthracene	17.42	178	1183	0.059	ng/ul#	84
15) Fluoranthene	19.35	202	13149	0.510	ng/ul	99
17) Pyrene	19.71	202	20070	0.736	ng/ul	99
18) Benzo(a)anthracene	21.45	228	13756	0.400	ng/ul	96
19) Chrysene	21.51	228	13345	0.502	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	32074m	1.098	ng/ul	
22) Benzo(k)fluoranthene	23.28	252	9362m	0.327	ng/ul	
23) Benzo(a)pyrene	23.91	252	11838	0.421	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.71	276	9766	0.303	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.73	278	2474	0.092	ng/ul#	73
26) Benzo(g,h,i)perylene	27.54	276	7305	0.270	ng/ul	98

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

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