

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097431.D
 Acq On : 7 Sep 2018 21:33
 Operator : SJ/JU
 Sample : J4688-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z32

Manual Integrations
 APPROVED

Sohil
 9/11/2018 2:43:25 PM

Quant Time: Sep 08 03:47:02 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 08 02:48:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1022	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5690	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	4012	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	10365m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10605m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	10529m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.71	152	2327	0.26	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	8270m	0.24	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.17	128	986	0.071	ng/ul#	88
5) 2-Methylnaphthalene	11.79	142	741	0.085	ng/ul	98
7) Acenaphthylene	13.73	152	948	0.065	ng/ul#	57
8) Acenaphthene	14.07	153	289	0.023	ng/ul#	83
9) Fluorene	15.07	166	495	0.034	ng/ul#	66
12) Phenanthrene	16.80	178	9146m	0.343	ng/ul	
13) Anthracene	16.89	178	2349m	0.105	ng/ul	
15) Fluoranthene	18.83	202	20852m	0.608	ng/ul	
17) Pyrene	19.20	202	22256m	0.780	ng/ul	
18) Benzo(a)anthracene	20.95	228	11517m	0.423	ng/ul	
19) Chrysene	21.00	228	9342m	0.298	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	43043m	1.527	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	11112m	0.353	ng/ul	
23) Benzo(a)pyrene	23.10	252	30276m	1.161	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.49	276	29475m	0.872	ng/ul	
25) Dibenzo(a,h)anthracene	25.50	278	7057m	0.246	ng/ul	
26) Benzo(g,h,i)perylene	26.19	276	31973m	1.109	ng/ul	

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

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