

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020818\
 Data File : BE095573.D
 Acq On : 8 Feb 2018 17:31
 Operator : SJ/JU
 Sample : J1315-20RE
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C83RE

Manual Integrations
 APPROVED

Sohil
 2/9/2018 4:30:53 PM

Quant Time: Feb 09 03:23:46 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 09 01:31:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1772	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5256	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	263780	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7894m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6188	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7449	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3296	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	9131	0.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	5825	0.396	ng/ul#	1
8) Acenaphthene	15.08	153	98585	0.100	ng/ul	93
9) Fluorene	16.03	166	115006m	0.113	ng/ul	
11) Pentachlorophenol	17.36	266	3500	2.289	ng/ul	99
12) Phenanthrene	17.75	178	18044m	0.680	ng/ul	
13) Anthracene	17.85	178	74890	3.082	ng/ul#	92
15) Fluoranthene	19.73	202	272425	7.925	ng/ul	80
17) Pyrene	20.08	202	214651m	10.338	ng/ul	
18) Benzo(a)anthracene	21.79	228	40726	2.105	ng/ul#	90
19) Chrysene	21.84	228	34507	1.718	ng/ul#	91
21) Benzo(b)fluoranthene	23.62	252	26718m	1.017	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	7252m	0.302	ng/ul	
23) Benzo(a)pyrene	24.32	252	15031m	0.641	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.23	276	5774	0.224	ng/ul#	75
25) Dibenzo(a,h)anthracene	27.24	278	1462	0.069	ng/ul#	15
26) Benzo(a,h,i)perylene	28.11	276	3889	0.178	ng/ul#	78

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

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