

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095424.D
 Acq On : 3 Feb 2018 7:17
 Operator : SJ/JU
 Sample : J1295-15DL 10X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C13DL

Manual Integrations
 APPROVED

Sohil
 2/5/2018 5:49:01 PM

Quant Time: Feb 05 01:32:06 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 03 05:26:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2154	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6461	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3917	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8576	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	8831	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	8725	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	447	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	1167	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.75	152	3478	0.184	ng/ul	99
9) Fluorene	16.02	166	622	0.041	ng/ul#	74
11) Pentachlorophenol	17.36	266	2208	1.329	ng/ul	97
12) Phenanthrene	17.75	178	958m	0.033	ng/ul	
13) Anthracene	17.84	178	2647	0.100	ng/ul	94
15) Fluoranthene	19.72	202	34669	0.928	ng/ul	84
17) Pyrene	20.07	202	56306	1.900	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	25895	0.938	ng/ul#	83
19) Chrysene	21.84	228	26510	0.925	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	35032m	1.138	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	11089m	0.394	ng/ul	
23) Benzo(a)pyrene	24.31	252	8295	0.302	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	8440	0.280	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.25	278	2074	0.084	ng/ul#	94
26) Benzo(g,h,i)perylene	28.11	276	6044	0.237	ng/ul	96

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

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