

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013018\
 Data File : BE095312.D
 Acq On : 31 Jan 2018 11:59
 Operator : SJ/JU
 Sample : J1244-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B42

Manual Integrations
 APPROVED

Sohil
 1/31/2018 8:29:02 PM

Quant Time: Jan 31 14:33:38 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 11:55:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2641	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	8251	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5729	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10811	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	11039	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	10310m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3336	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	8763	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	108699	4.705	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	53084	3.441	ng/ul	100
7) Acenaphthylene	14.74	152	3102	0.113	ng/ul#	92
8) Acenaphthene	15.08	153	48220	2.255	ng/ul	94
9) Fluorene	16.03	166	28712	1.302	ng/ul#	78
11) Pentachlorophenol	17.36	266	92393	44.124	ng/ul	95
12) Phenanthrene	17.75	178	165385	4.549	ng/ul#	86
13) Anthracene	17.85	178	23432	0.704	ng/ul#	93
15) Fluoranthene	19.73	202	422740	8.979	ng/ul	81
17) Pyrene	20.08	202	316738	8.551	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	62396	1.808	ng/ul#	86
19) Chrysene	21.84	228	48983	1.367	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	32712m	0.899	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	12785m	0.384	ng/ul	
23) Benzo(a)pyrene	24.32	252	14072m	0.433	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	6907m	0.194	ng/ul	
25) Dibenzo(a,h)anthracene	27.26	278	2060m	0.071	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	5626m	0.187	ng/ul	

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

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