

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095401.D
 Acq On : 2 Feb 2018 17:36
 Operator : SJ/JU
 Sample : J1243-06DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B29DL

Manual Integrations
 APPROVED

Sohil
 2/5/2018 5:48:00 PM

Quant Time: Feb 03 04:47:51 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 04:31:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1268m	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5263	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3156	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7022	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7261	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7375	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	569	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	1547	0.06	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.31	128	874	0.059	ng/ul#	76
5) 2-Methylnaphthalene	12.89	142	728	0.074	ng/ul	99
7) Acenaphthylene	14.74	152	3188	0.210	ng/ul	99
8) Acenaphthene	15.08	153	1069	0.091	ng/ul#	95
9) Fluorene	16.02	166	1631	0.134	ng/ul	85
11) Pentachlorophenol	17.36	266	25436	18.702	ng/ul	97
12) Phenanthrene	17.75	178	20389	0.863	ng/ul#	90
13) Anthracene	17.85	178	30455	1.409	ng/ul#	92
15) Fluoranthene	19.72	202	95108	3.110	ng/ul	82
17) Pyrene	20.08	202	86140	3.535	ng/ul#	77
18) Benzo(a)anthracene	21.79	228	19039	0.839	ng/ul#	88
19) Chrysene	21.84	228	19705	0.836	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	23911m	0.919	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	6853m	0.288	ng/ul	
23) Benzo(a)pyrene	24.31	252	10600	0.456	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	27.23	276	6811	0.267	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.23	278	1615	0.077	ng/ul#	76
26) Benzo(g,h,i)perylene	28.12	276	5037	0.233	ng/ul	96

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

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