

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097875.D
 Acq On : 12 Oct 2018 2:04
 Operator : SJ/JU
 Sample : J5183-01DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC16DL

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:35:05 PM

Quant Time: Oct 12 05:19:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 04:56:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2774	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	14046	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	9550	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.31	188	23257	0.40	ng/ul	0.00
16) Chrysene-d12	21.49	240	24203	0.40	ng/ul	0.00
20) Perylene-d12	24.06	264	25448	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1473	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.34	212	4718	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	778	0.023	ng/ul#	69
5) 2-Methylnaphthalene	12.37	142	571	0.023	ng/ul	98
8) Acenaphthene	14.62	153	786	0.028	ng/ul	96
12) Phenanthrene	17.35	178	17715	0.322	ng/ul	98
13) Anthracene	17.44	178	3097	0.061	ng/ul#	89
15) Fluoranthene	19.37	202	52079	0.737	ng/ul	97
17) Pyrene	19.73	202	62454	0.866	ng/ul	99
18) Benzo(a)anthracene	21.48	228	31019	0.424	ng/ul#	92
19) Chrysene	21.53	228	50687	0.771	ng/ul#	92
21) Benzo(b)fluoranthene	23.27	252	68443m	1.006	ng/ul	
22) Benzo(k)fluoranthene	23.32	252	18281m	0.251	ng/ul	
23) Benzo(a)pyrene	23.95	252	27628	0.414	ng/ul#	68
24) Indeno(1,2,3-cd)pyrene	26.76	276	26793	0.353	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.79	278	6875	0.110	ng/ul#	7
26) Benzo(g,h,i)perylene	27.60	276	30330	0.459	ng/ul#	83

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

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