

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098090.D
 Acq On : 26 Oct 2018 16:02
 Operator : SJ/JU
 Sample : J5598-05DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE8DL

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:23 PM

Quant Time: Oct 27 00:25:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 26 23:15:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	1003	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	5059	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.54	164	3529	0.40	ng/ul	0.01
10) Phenanthrene-d10	17.30	188	8362	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11351	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	9856	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	372	0.04	ng/ul	0.01
14) Fluoranthene-d10	19.32	212	1095	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	1347	0.111	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	278	0.030	ng/ul	95
7) Acenaphthylene	14.27	152	1084	0.078	ng/ul#	89
8) Acenaphthene	14.61	153	358	0.034	ng/ul#	88
12) Phenanthrene	17.34	178	3274	0.170	ng/ul	95
13) Anthracene	17.43	178	1786	0.094	ng/ul#	86
15) Fluoranthene	19.36	202	26831	1.090	ng/ul	99
17) Pyrene	19.72	202	32477	1.225	ng/ul	99
18) Benzo(a)anthracene	21.46	228	23213	0.696	ng/ul	97
19) Chrysene	21.52	228	19875	0.770	ng/ul	98
21) Benzo(b)fluoranthene	23.25	252	44136m	1.616	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	14753m	0.552	ng/ul	
23) Benzo(a)pyrene	23.91	252	18212	0.693	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.72	276	14800	0.492	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.74	278	4067	0.162	ng/ul#	86
26) Benzo(g,h,i)perylene	27.55	276	10363	0.409	ng/ul	96

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

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