

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102114\
 Data File : BE088151.D
 Acq On : 21 Oct 2014 20:14
 Operator : TP/JJ
 Sample : F4351-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 YS19-SW08P-1014

Quant Time: Oct 22 05:00:25 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 21 02:11:14 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.14	152	23816	5.00	ng	0.00
3) Naphthalene-d8	8.70	136	92228	5.00	ng	0.00
7) Acenaphthene-d10	10.85	164	43844	5.00	ng	0.00
12) Phenanthrene-d10	12.96	188	81914	5.00	ng	0.00
16) Chrysene-d12	18.38	240	51139	5.00	ng	0.00
22) Perylene-d12	21.44	264	40339	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.83	82	46131	6.94	ng	0.00
8) 2-Fluorobiphenyl	10.03	172	83560	6.90	ng	0.00
18) Terphenyl-d14	16.15	244	81213	9.87	ng	0.00
Target Compounds						Qvalue
14) Anthracene	13.09	178	5301	0.08	ng	96
15) Fluoranthene	15.25	202	1860	0.12	ng	100
17) Pyrene	15.70	202	2002	0.14	ng	99
19) Benzo(a)anthracene	18.36	228	5449	0.12	ng	# 90
20) Chrysene	18.44	228	8166	0.18	ng	# 91
21) Indeno(1,2,3-cd)pyrene	23.63	276	3198	0.06	ng	98
23) Benzo(b)fluoranthene	20.69	252	1737	0.18	ng	# 82
24) Benzo(k)fluoranthene	20.74	252	1076	0.10	ng	# 74
27) Benzo(g,h,i)perylene	24.26	276	3135	0.08	ng	# 68

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

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