

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088461.D
 Acq On : 1 Nov 2014 20:57
 Operator : TP/JJ
 Sample : F4368-17RE
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 YS19-SB44-1014RE

Quant Time: Nov 02 04:57:35 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE103114.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 01 00:42:07 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.09	152	31290	5.00	ng	0.00
3) Naphthalene-d8	8.65	136	111361	5.00	ng	0.00
7) Acenaphthene-d10	10.79	164	51650	5.00	ng	0.00
12) Phenanthrene-d10	12.91	188	107776	5.00	ng	0.00
16) Chrysene-d12	18.33	240	66617	5.00	ng	0.01
22) Perylene-d12	21.39	264	64589	5.00	ng	0.02
System Monitoring Compounds						
4) Nitrobenzene-d5	7.79	82	498391	61.03	ng	0.02
8) 2-Fluorobiphenyl	9.99	172	569098	39.09	ng	0.00
18) Terphenyl-d14	16.11	244	484229	39.57	ng	0.03
Target Compounds						Qvalue
13) Phenanthrene	12.95	178	10924	0.11	ng	# 55
15) Fluoranthene	15.18	202	7055	0.44	ng	100
17) Pyrene	15.64	202	6766	0.31	ng	98
19) Benzo(a)anthracene	18.31	228	12450	0.20	ng	# 88
20) Chrysene	18.38	228	22939	0.37	ng	# 93
21) Indeno(1,2,3-cd)pyrene	23.57	276	13216	0.23	ng	# 84
23) Benzo(b)fluoranthene	20.63	252	8802	0.62	ng	# 80
24) Benzo(k)fluoranthene	20.69	252	2811	0.18	ng	# 58
25) Benzo(a)pyrene	21.27	252	3517	0.26	ng	# 61
27) Benzo(g,h,i)perylene	24.20	276	13745	0.24	ng	# 77

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

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