

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088468.D
 Acq On : 2 Nov 2014 00:50
 Operator : TP/JJ
 Sample : F4368-03RE
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 YS19-SS33-1014RE

Quant Time: Nov 02 05:29:15 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	22930	5.00	ng	0.00
3) Naphthalene-d8	8.65	136	80016	5.00	ng	0.00
7) Acenaphthene-d10	10.80	164	36775	5.00	ng	0.01
12) Phenanthrene-d10	12.91	188	74811	5.00	ng	0.00
16) Chrysene-d12	18.33	240	47604	5.00	ng	0.02
22) Perylene-d12	21.41	264	44936	5.00	ng	0.03
System Monitoring Compounds						
4) Nitrobenzene-d5	7.79	82	523387	89.19	ng	0.02
8) 2-Fluorobiphenyl	9.99	172	470744	45.41	ng	0.00
18) Terphenyl-d14	16.10	244	316785	36.22	ng	0.03
Target Compounds						Qvalue
13) Phenanthrene	12.95	178	15779	0.23	ng	# 76
14) Anthracene	13.03	178	2284	0.04	ng	# 66
15) Fluoranthene	15.19	202	5023	0.46	ng	98
17) Pyrene	15.65	202	4152	0.27	ng	97
19) Benzo(a)anthracene	18.31	228	7538	0.17	ng	# 83
20) Chrysene	18.38	228	9371	0.21	ng	# 86
21) Indeno(1,2,3-cd)pyrene	23.59	276	5175	0.13	ng	# 52

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

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