

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080416\
 Data File : BE092193.D
 Acq On : 4 Aug 2016 21:47
 Operator : UM/SJ
 Sample : H4269-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 GW-52-11.1-15.0

Quant Time: Aug 05 03:56:32 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-PAH-BE080216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 16:59:37 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	10631	5.00	ng	0.00
6) Naphthalene-d8	10.98	136	47900	5.00	ng	-0.02
10) Acenaphthene-d10	14.83	164	26276	5.00	ng	-0.02
16) Phenanthrene-d10	17.56	188	62412	5.00	ng	-0.02
22) Chrysene-d12	21.75	240	79325	5.00	ng	0.00
28) Perylene-d12	24.14	264	78653	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.69	112	11366	4.16	ng	0.00
5) Phenol-d6	7.35	99	9443	2.60	ng	0.00
7) Nitrobenzene-d5	9.34	82	12417	3.29	ng	-0.01
11) 2,4,6-Tribromophenol	16.33	330	6814	7.99	ng	0.00
12) 2-Fluorobiphenyl	13.46	172	28055	3.37	ng	0.00
24) Terphenyl-d14	20.19	244	46922	4.12	ng	-0.02
Target Compounds						
23) Pyrene	19.99	202	1662	0.02	ng	# 79
27) Indeno(1,2,3-cd)pyrene	26.65	276	2018	0.02	ng	97
33) Benzo(a,h,i)perylene	27.42	276	3026	0.04	ng	# 63

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

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