

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
 Data File : BF080156.D
 Acq On : 25 Jun 2015 5:02
 Operator : TP/IZ
 Sample : G2699-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LAR9819

Quant Time: Jun 25 05:44:44 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 00:52:28 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.32	152	33838	20.00	ng	0.00
21) Naphthalene-d8	8.61	136	130578	20.00	ng	0.00
38) Acenaphthene-d10	10.38	164	66830	20.00	ng	0.00
63) Phenanthrene-d10	11.88	188	128818	20.00	ng	-0.01
75) Chrysene-d12	14.78	240	158801	20.00	ng	-0.15
86) Perylene-d12	16.62	264	162166	20.00	ng	-0.22
System Monitoring Compounds						
5) 2-Fluorophenol	5.92	112	292482	153.01	ng	0.00
7) Phenol-d6	6.92	99	367764	144.58	ng	0.00
23) Nitrobenzene-d5	7.88	82	215851	96.23	ng	0.00
41) 2,4,6-Tribromophenol	11.18	330	97333	148.94	ng	0.01
44) 2-Fluorobiphenyl	9.68	172	439241	93.73	ng	0.00
78) Terphenyl-d14	13.54	244	570509	83.90	ng	-0.07
Target Compounds						Qvalue
49) Dimethylphthalate	10.07	163	50638	10.08	ng	# 98
70) Phenanthrene	11.90	178	17762	2.28	ng	96
74) Fluoranthene	13.14	202	52601	6.33	ng	91
77) Pyrene	13.40	202	52863	5.41	ng	97
80) Benzo(a)anthracene	14.77	228	44049m	4.55	ng	
82) Chrysene	14.81	228	42193m	4.73	ng	
85) Indeno(1,2,3-cd)pyrene	18.45	276	46510	4.13	ng	# 82
87) Benzo(b)fluoranthene	16.08	252	75862m	7.73	ng	
89) Benzo(a)pyrene	16.54	252	53659m	5.75	ng	
91) Benzo(g,h,i)perylene	19.00	276	49176	5.10	ng	# 78

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

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