

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
 Data File : BF088207.D
 Acq On : 21 Jun 2016 8:59
 Operator : UM/SJ
 Sample : H3570-20RE
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-22-COMPRES

Manual Integrations
 APPROVED

sohil
 6/21/2016 4:22:38 PM

Quant Time: Jun 21 12:54:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 17:39:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	149319	20.00	ng	-0.01
21) Naphthalene-d8	7.95	136	605261	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	281542	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	448792	20.00	ng	0.00
75) Chrysene-d12	13.81	240	328630	20.00	ng	0.00
86) Perylene-d12	15.18	264	338984	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	296401	33.93	ng	0.00
7) Phenol-d6	6.32	99	435562	41.15	ng	-0.01
23) Nitrobenzene-d5	7.23	82	241322	23.28	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	182175	61.28	ng	0.00
44) 2-Fluorobiphenyl	9.03	172	519503	25.08	ng	-0.01
78) Terphenyl-d14	12.75	244	637747	44.16	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.43	163	266221	13.05	ng	99
77) Pyrene	12.61	202	70688	2.90	ng	98
80) Benzo(a)anthracene	13.79	228	55448	2.96	ng	99
82) Chrysene	13.83	228	40798	2.32	ng	98
85) Indeno(1,2,3-cd)pyrene	16.47	276	50906	3.11	ng	# 100
87) Benzo(b)fluoranthene	14.80	252	89334m	3.78	ng	
89) Benzo(a)pyrene	15.13	252	62164	3.25	ng	97
91) Benzo(g,h,i)perylene	16.86	276	53145	2.91	ng	98

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

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