

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071116\
 Data File : BF088909.D
 Acq On : 11 Jul 2016 19:54
 Operator : UM/SJ
 Sample : H3915-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P6-B1-0-11-C

Manual Integrations

sohil
 7/12/2016 7:35:51 PM

Quant Time: Jul 12 02:29:35 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 11 18:13:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	67050	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	270731	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	106983	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	178685	20.00	ng	0.00
75) Chrysene-d12	13.84	240	137205	20.00	ng	-0.01
86) Perylene-d12	15.21	264	124536	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	594978	132.99	ng	0.02
7) Phenol-d6	6.36	99	831499	143.84	ng	0.01
23) Nitrobenzene-d5	7.28	82	449998	87.10	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	148370	149.35	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	623943	92.12	ng	0.00
78) Terphenyl-d14	12.81	244	466889	75.79	ng	0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.47	163	157618	20.68	ng	99
70) Phenanthrene	11.24	178	56597	5.79	ng	97
74) Fluoranthene	12.42	202	118641	11.98	ng	99
77) Pyrene	12.65	202	99191	8.53	ng	97
80) Benzo(a)anthracene	13.83	228	63199	7.15	ng	94
82) Chrysene	13.86	228	41439m	4.74	ng	
85) Indeno(1,2,3-cd)pyrene	16.49	276	18925	2.81	ng	# 100
87) Benzo(b)fluoranthene	14.83	252	49655m	6.36	ng	
88) Benzo(k)fluoranthene	14.85	252	22959m	3.38	ng	
89) Benzo(a)pyrene	15.15	252	33482	5.06	ng	99
91) Benzo(g,h,i)perylene	16.88	276	15759	2.76	ng	96

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

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