

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092616\
 Data File : BF090237.D
 Acq On : 26 Sep 2016 22:29
 Operator : UM/SJ
 Sample : H4898-01RE
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CAM-SELECTFILLRE

Manual Integrations
 APPROVED

sohil
 9/27/2016 1:06:37 PM

Quant Time: Sep 27 01:46:04 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 26 15:58:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.48	152	161391	20.00	ng	0.00
21) Naphthalene-d8	7.77	136	617858	20.00	ng	0.00
38) Acenaphthene-d10	9.51	164	287674	20.00	ng	0.00
63) Phenanthrene-d10	10.98	188	400554	20.00	ng	0.00
75) Chrysene-d12	13.59	240	287054	20.00	ng	-0.01
86) Perylene-d12	14.91	264	224194	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.04	112	228215	23.05	ng	0.01
7) Phenol-d6	6.14	99	876472	71.68	ng	0.00
23) Nitrobenzene-d5	7.05	82	678410	63.65	ng	-0.01
41) 2,4,6-Tribromophenol	10.30	330	2429	0.98	ng	0.00
44) 2-Fluorobiphenyl	8.86	172	1299135	88.66	ng	0.00
78) Terphenyl-d14	12.56	244	780250	69.01	ng	0.00
Target Compounds						
10) Phenol	6.15	94	91788	6.35	ng	Qvalue # 69
49) Dimethylphthalate	9.24	163	360757	19.69	ng	99
70) Phenanthrene	11.01	178	120117	6.41	ng	98
74) Fluoranthene	12.17	202	131451	6.54	ng	97
77) Pyrene	12.40	202	116745	5.94	ng	97
80) Benzo(a)anthracene	13.58	228	61748	4.00	ng	# 91
82) Chrysene	13.61	228	45759m	3.04	ng	
87) Benzo(b)fluoranthene	14.57	252	47779m	3.85	ng	
88) Benzo(k)fluoranthene	14.58	252	23624m	2.03	ng	
89) Benzo(a)pyrene	14.87	252	35286	3.02	ng	# 62
91) Benzo(g,h,i)perylene	16.40	276	20272	2.27	ng	# 81

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

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