

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111116\
 Data File : BF091156.D
 Acq On : 11 Nov 2016 22:35
 Operator : UM/SJ
 Sample : H5628-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-COMP

Manual Integrations
 APPROVED

sohil
 11/14/2016 6:36:46 PM

Quant Time: Nov 12 00:48:43 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 12:59:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	150085	20.00	ng	-0.01
21) Naphthalene-d8	7.93	136	659737	20.00	ng	-0.02
38) Acenaphthene-d10	9.68	164	317757	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	493127	20.00	ng	-0.01
75) Chrysene-d12	13.78	240	367263	20.00	ng	-0.03
86) Perylene-d12	15.15	264	274193	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	1083480	119.23	ng	-0.01
7) Phenol-d6	6.30	99	1483241	120.25	ng	-0.01
23) Nitrobenzene-d5	7.21	82	879751	72.74	ng	-0.03
41) 2,4,6-Tribromophenol	10.46	330	308920	107.54	ng	-0.02
44) 2-Fluorobiphenyl	9.01	172	1413157	76.57	ng	-0.02
78) Terphenyl-d14	12.74	244	1189813	80.84	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	69278	5.07	ng	Qvalue # 60
49) Dimethylphthalate	9.41	163	95981	4.61	ng	98
70) Phenanthrene	11.18	178	76133	2.90	ng	97
74) Fluoranthene	12.36	202	179794	6.61	ng	93
77) Pyrene	12.59	202	178839	7.44	ng	96
80) Benzo(a)anthracene	13.78	228	78813m	3.78	ng	
82) Chrysene	13.81	228	63517m	3.09	ng	
87) Benzo(b)fluoranthene	14.79	252	61201m	3.29	ng	
89) Benzo(a)pyrene	15.11	252	46159	2.85	ng	# 97
91) Benzo(g,h,i)perylene	16.81	276	29791	2.35	ng	97

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

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