

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072316\
 Data File : BF089264.D
 Acq On : 24 Jul 2016 7:05
 Operator : UM/SJ
 Sample : H4151-10
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-05-COMP

Manual Integrations

sohil
 7/25/2016 6:43:27 PM

Quant Time: Jul 25 05:19:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 03:29:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.58	152	48937	20.00	ng	0.00
21) Naphthalene-d8	7.86	136	215276	20.00	ng	-0.01
38) Acenaphthene-d10	9.61	164	90542	20.00	ng	0.00
63) Phenanthrene-d10	11.08	188	145533	20.00	ng	0.00
75) Chrysene-d12	13.70	240	98686	20.00	ng	-0.01
86) Perylene-d12	15.05	264	93530	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	368173	114.42	ng	0.01
7) Phenol-d6	6.24	99	507685	122.35	ng	0.00
23) Nitrobenzene-d5	7.15	82	306007	75.53	ng	0.00
41) 2,4,6-Tribromophenol	10.40	330	77986	95.67	ng	0.00
44) 2-Fluorobiphenyl	8.95	172	517656	78.77	ng	0.00
78) Terphenyl-d14	12.66	244	289843	63.70	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.34	163	75579	10.88	ng	Qvalue 96
74) Fluoranthene	12.29	202	30559	3.64	ng	96
77) Pyrene	12.51	202	27901	3.42	ng	97
80) Benzo(a)anthracene	13.69	228	17886	2.85	ng	# 90
82) Chrysene	13.73	228	13517m	2.25	ng	
85) Indeno(1,2,3-cd)pyrene	16.26	276	9665	2.22	ng	95
87) Benzo(b)fluoranthene	14.70	252	19324m	2.85	ng	
88) Benzo(k)fluoranthene	14.70	252	12626m	2.61	ng	
89) Benzo(a)pyrene	14.99	252	16708	3.18	ng	96
91) Benzo(g,h,i)perylene	16.63	276	8489	2.03	ng	96

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

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