

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072316\
 Data File : BF089262.D
 Acq On : 24 Jul 2016 6:05
 Operator : UM/SJ
 Sample : H4151-06
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-07-COMP

Manual Integrations

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

Quant Time: Jul 25 05:15:00 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	46788	20.00	ng	0.00
21) Naphthalene-d8	7.86	136	214531	20.00	ng	-0.01
38) Acenaphthene-d10	9.61	164	87682	20.00	ng	0.00
63) Phenanthrene-d10	11.08	188	135019	20.00	ng	0.00
75) Chrysene-d12	13.70	240	96911	20.00	ng	-0.01
86) Perylene-d12	15.05	264	97073	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	359194	116.76	ng	0.01
7) Phenol-d6	6.24	99	496623	125.18	ng	0.00
23) Nitrobenzene-d5	7.15	82	300895	74.52	ng	0.00
41) 2,4,6-Tribromophenol	10.40	330	76621	97.06	ng	0.00
44) 2-Fluorobiphenyl	8.95	172	523027	82.19	ng	0.00
78) Terphenyl-d14	12.66	244	286537	64.13	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.34	163	59999	8.92	ng	Qvalue # 96
70) Phenanthrene	11.11	178	23076	3.12	ng	99
74) Fluoranthene	12.29	202	48386	6.22	ng	96
77) Pyrene	12.50	202	40798m	5.10	ng	
80) Benzo(a)anthracene	13.69	228	24011	3.89	ng	93
82) Chrysene	13.73	228	20214	3.43	ng	97
85) Indeno(1,2,3-cd)pyrene	16.26	276	10658	2.49	ng	# 90
87) Benzo(b)fluoranthene	14.69	252	22625m	3.22	ng	
88) Benzo(k)fluoranthene	14.70	252	14471m	2.88	ng	
89) Benzo(a)pyrene	14.99	252	17486	3.20	ng	95

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

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