

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
 Data File : BF085363.D
 Acq On : 11 Mar 2016 5:34
 Operator : UM/SJ
 Sample : H1731-08
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-21-COMP

Manual Integrations
 APPROVED

Sohil
 3/11/2016 4:58:49 PM

Quant Time: Mar 11 06:11:02 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 02:49:26 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	159050	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	564308	20.00	ng	-0.01
38) Acenaphthene-d10	9.98	164	252141	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	410312	20.00	ng	0.00
75) Chrysene-d12	14.09	240	268793	20.00	ng	-0.01
86) Perylene-d12	15.56	264	270100	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1198096	126.36	ng	0.00
7) Phenol-d6	6.57	99	1775236	142.90	ng	0.00
23) Nitrobenzene-d5	7.50	82	938915	89.03	ng	-0.01
41) 2,4,6-Tribromophenol	10.77	330	289993	134.41	ng	0.00
44) 2-Fluorobiphenyl	9.30	172	1596366	96.29	ng	0.00
78) Terphenyl-d14	13.05	244	1077751	93.08	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.69	163	265229	13.70	ng	99
74) Fluoranthene	12.68	202	70189	3.25	ng	94
77) Pyrene	12.90	202	66334	3.28	ng	96
80) Benzo(a)anthracene	14.08	228	41188	2.56	ng	94
82) Chrysene	14.13	228	32503m	2.19	ng	
85) Indeno(1,2,3-cd)pyrene	17.01	276	24206	2.09	ng	92
87) Benzo(b)fluoranthene	15.13	252	51534m	2.86	ng	
89) Benzo(a)pyrene	15.50	252	37498	2.51	ng	97

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

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