

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030216\
 Data File : BF085118.D
 Acq On : 2 Mar 2016 14:40
 Operator : SJ/IZ
 Sample : H1631-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-43(6-8)

Manual Integrations
 APPROVED

UMANGI
 3/3/2016 2:05:29 PM

Quant Time: Mar 03 12:39:44 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 02 13:25:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	146009	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	631541	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	240436	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	342219	20.00	ng	0.00
75) Chrysene-d12	14.18	240	308207	20.00	ng	0.03
86) Perylene-d12	15.69	264	283014	20.00	ng	0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	669885	74.48	ng	0.00
7) Phenol-d6	6.60	99	886175	75.16	ng	-0.02
23) Nitrobenzene-d5	7.54	82	507401	42.74	ng	-0.02
41) 2,4,6-Tribromophenol	10.82	330	183805	76.07	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	787513	48.24	ng	0.00
78) Terphenyl-d14	13.11	244	538744	41.01	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.74	163	184160	10.33	ng	# 93
51) Acenaphthene	10.07	154	91584	6.31	ng	95
77) Pyrene	12.97	202	596636	28.36	ng	95
85) Indeno(1,2,3-cd)pyrene	17.20	276	89833	6.50	ng	# 83
89) Benzo(a)pyrene	15.56	252	226325m	15.62	ng	
91) Benzo(g,h,i)perylene	17.65	276	72338	5.45	ng	# 63

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

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