

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030316\
 Data File : BF085198.D
 Acq On : 4 Mar 2016 11:16
 Operator : SJ/IZ
 Sample : H1631-02RE
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 3/4/2016 4:36:56 PM

Quant Time: Mar 08 10:28:35 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:54:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	130407	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	493412	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	168438	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	258318	20.00	ng	0.01
75) Chrysene-d12	14.17	240	230652	20.00	ng	0.02
86) Perylene-d12	15.69	264	197392	20.00	ng	0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	568792	73.16	ng	0.00
7) Phenol-d6	6.60	99	694536	68.19	ng	-0.01
23) Nitrobenzene-d5	7.54	82	433927	47.06	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	102593	71.18	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	606026	54.72	ng	-0.01
78) Terphenyl-d14	13.11	244	426864	42.96	ng	0.02
Target Compounds						
49) Dimethylphthalate	9.74	163	139995	10.83	ng	# 79
51) Acenaphthene	10.06	154	69966	6.75	ng	94
77) Pyrene	12.97	202	517428	29.81	ng	96
85) Indeno(1,2,3-cd)pyrene	17.20	276	60493	6.08	ng	# 84
89) Benzo(a)pyrene	15.56	252	163725m	15.02	ng	
91) Benzo(g,h,i)perylene	17.66	276	49548	5.29	ng	# 55

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

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