

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
 Data File : BF087827.D
 Acq On : 8 Jun 2016 23:45
 Operator : UM/SJ
 Sample : H3399-07 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-28

Manual Integrations
 APPROVED

sohil
 6/9/2016 8:04:05 PM

Quant Time: Jun 09 13:19:27 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	74056	20.00	ng	-0.02
21) Naphthalene-d8	8.11	136	316900	20.00	ng	-0.01
38) Acenaphthene-d10	9.87	164	132462	20.00	ng	-0.01
63) Phenanthrene-d10	11.35	188	225798	20.00	ng	-0.01
75) Chrysene-d12	13.99	240	201358	20.00	ng	0.00
86) Perylene-d12	15.42	264	157483	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	239041	55.18	ng	-0.01
7) Phenol-d6	6.46	99	297390	56.65	ng	-0.02
23) Nitrobenzene-d5	7.38	82	141011	25.98	ng	-0.03
41) 2,4,6-Tribromophenol	10.65	330	26135	18.69	ng	-0.02
44) 2-Fluorobiphenyl	9.19	172	310982	33.53	ng	-0.02
78) Terphenyl-d14	12.94	244	257756	29.13	ng	-0.01
Target Compounds						Qvalue
15) Benzyl Alcohol	6.97	79	48232	11.74	ng	91
49) Dimethylphthalate	9.59	163	29884	3.11	ng	100
70) Phenanthrene	11.38	178	198762	16.78	ng	98
71) Anthracene	11.43	178	65588	5.49	ng	97
74) Fluoranthene	12.56	202	299824	23.98	ng	95
77) Pyrene	12.79	202	266454	17.83	ng	98
80) Benzo(a)anthracene	13.98	228	159347	13.89	ng	98
82) Chrysene	14.01	228	107321m	9.94	ng	
85) Indeno(1,2,3-cd)pyrene	16.80	276	46284	4.61	ng	# 100
87) Benzo(b)fluoranthene	15.01	252	149061m	13.58	ng	
88) Benzo(k)fluoranthene	15.03	252	26637m	3.22	ng	
89) Benzo(a)pyrene	15.36	252	86365	9.73	ng	# 96
91) Benzo(a,h,i)perylene	17.22	276	45202	5.33	ng	96

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

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