

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080116\
 Data File : BF089535.D
 Acq On : 2 Aug 2016 3:25
 Operator : UM/SJ
 Sample : H4288-12
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-66-0.5-1.5

Manual Integrations
 APPROVED

sohil
 8/2/2016 5:06:46 PM

Quant Time: Aug 02 04:01:53 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 02:21:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.48	152	34165	20.00	ng	0.00
21) Naphthalene-d8	7.77	136	115689	20.00	ng	0.00
38) Acenaphthene-d10	9.51	164	48779	20.00	ng	-0.01
63) Phenanthrene-d10	10.98	188	81958	20.00	ng	0.00
75) Chrysene-d12	13.60	240	92120	20.00	ng	-0.01
86) Perylene-d12	14.93	264	78110	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.05	112	226933	106.06	ng	0.01
7) Phenol-d6	6.15	99	277953	98.30	ng	0.00
23) Nitrobenzene-d5	7.05	82	168958	86.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.30	330	37813	93.60	ng	-0.01
44) 2-Fluorobiphenyl	8.84	172	266589	80.21	ng	-0.01
78) Terphenyl-d14	12.56	244	201873	51.29	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.24	163	44259	11.92	ng	99
70) Phenanthrene	11.01	178	68555	16.22	ng	98
71) Anthracene	11.05	178	24790	5.27	ng	98
73) Di-n-butylphthalate	11.57	149	20500	3.86	ng	# 97
74) Fluoranthene	12.18	202	155113	34.89	ng	99
77) Pyrene	12.40	202	143615	20.57	ng	97
80) Benzo(a)anthracene	13.59	228	90842	17.47	ng	97
82) Chrysene	13.62	228	69917	12.63	ng	98
85) Indeno(1,2,3-cd)pyrene	16.09	276	36733	9.34	ng	# 85
87) Benzo(b)fluoranthene	14.59	252	109366m	22.55	ng	
88) Benzo(k)fluoranthene	14.61	252	23087m	5.18	ng	
89) Benzo(a)pyrene	14.88	252	66406	15.26	ng	97
90) Dibenzo(a,h)anthracene	16.09	278	13788	4.02	ng	# 79
91) Benzo(g,h,i)perylene	16.43	276	36230	10.75	ng	99

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

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