

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
 Data File : BF089498.D
 Acq On : 1 Aug 2016 6:30
 Operator : UM/SJ
 Sample : H4215-08
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11-COMP

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:16:56 PM

Quant Time: Aug 01 07:33:18 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 04:11:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	37764	20.00	ng	-0.01
21) Naphthalene-d8	7.78	136	134845	20.00	ng	-0.01
38) Acenaphthene-d10	9.52	164	56389	20.00	ng	-0.01
63) Phenanthrene-d10	10.99	188	111327	20.00	ng	-0.01
75) Chrysene-d12	13.62	240	88512	20.00	ng	0.00
86) Perylene-d12	14.96	264	61597	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.07	112	255871	108.19	ng	0.01
7) Phenol-d6	6.16	99	329611	105.46	ng	-0.01
23) Nitrobenzene-d5	7.06	82	187541	82.09	ng	-0.02
41) 2,4,6-Tribromophenol	10.31	330	48018	102.82	ng	-0.01
44) 2-Fluorobiphenyl	8.86	172	275385	71.67	ng	-0.01
78) Terphenyl-d14	12.58	244	256565	67.84	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	7.80	128	16160	2.42	ng	99
48) Acenaphthylene	9.38	152	64231	11.31	ng	98
49) Dimethylphthalate	9.26	163	47543	11.08	ng	97
51) Acenaphthene	9.55	154	6857	2.07	ng	99
57) Fluorene	10.07	166	25254	6.40	ng	98
70) Phenanthrene	11.02	178	127518	22.21	ng	98
71) Anthracene	11.07	178	61911	9.70	ng	100
72) Carbazole	11.23	167	17576	3.27	ng	# 93
74) Fluoranthene	12.21	202	279817	46.33	ng	98
77) Pyrene	12.42	202	236307	35.23	ng	98
80) Benzo(a)anthracene	13.61	228	135706	27.17	ng	# 86
82) Chrysene	13.65	228	98470m	18.52	ng	
83) Bis(2-ethylhexyl)phthalate	13.63	149	38227	8.67	ng	# 94
85) Indeno(1,2,3-cd)pyrene	16.15	276	54832	14.50	ng	98
87) Benzo(b)fluoranthene	14.62	252	125698m	32.86	ng	
88) Benzo(k)fluoranthene	14.63	252	42968m	12.22	ng	
89) Benzo(a)pyrene	14.91	252	81031	23.61	ng	96
90) Dibenzo(a,h)anthracene	16.15	278	15316	5.67	ng	# 80
91) Benzo(a,h,i)perylene	16.50	276	55282	20.80	ng	97

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

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