

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060616\
 Data File : BF087758.D
 Acq On : 6 Jun 2016 18:15
 Operator : UM/SJ
 Sample : H3382-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-02-COMP

Manual Integrations
 APPROVED

sohil
 6/7/2016 7:22:15 PM

Quant Time: Jun 07 01:39:04 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 04 11:07:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	104136	20.00	ng	-0.02
21) Naphthalene-d8	8.14	136	432247	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	167246	20.00	ng	-0.02
63) Phenanthrene-d10	11.37	188	279471	20.00	ng	-0.01
75) Chrysene-d12	14.00	240	210791	20.00	ng	-0.02
86) Perylene-d12	15.43	264	172394	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	649684	100.84	ng	0.00
7) Phenol-d6	6.49	99	938274	116.12	ng	0.00
23) Nitrobenzene-d5	7.42	82	547539	73.29	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	152703	90.96	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	898504	81.31	ng	-0.02
78) Terphenyl-d14	12.96	244	651201	75.38	ng	-0.01
Target Compounds						Qvalue
48) Acenaphthylene	9.75	152	64666	3.84	ng	99
49) Dimethylphthalate	9.61	163	69459	5.71	ng	98
70) Phenanthrene	11.39	178	303985	20.06	ng	99
71) Anthracene	11.44	178	61363	4.04	ng	99
72) Carbazole	11.60	167	30305	2.11	ng	98
74) Fluoranthene	12.58	202	561666	37.67	ng	96
77) Pyrene	12.81	202	427683	28.49	ng	98
80) Benzo(a)anthracene	13.99	228	283335m	23.29	ng	
82) Chrysene	14.02	228	196698	16.24	ng	97
85) Indeno(1,2,3-cd)pyrene	16.81	276	99488	9.94	ng	# 100
87) Benzo(b)fluoranthene	15.02	252	257972m	23.00	ng	
88) Benzo(k)fluoranthene	15.03	252	115802m	12.41	ng	
89) Benzo(a)pyrene	15.36	252	160816	17.06	ng	99
90) Dibenzo(a,h)anthracene	16.82	278	26517	3.31	ng	94
91) Benzo(a,h,i)perylene	17.22	276	85792	10.60	ng	99

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

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