

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010317\
 Data File : BF092068.D
 Acq On : 3 Jan 2017 18:46
 Operator : UM/SJ
 Sample : H6325-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LAR1020

Manual Integrations
 APPROVED

Sohil
 1/4/2017 2:05:48 PM

Quant Time: Jan 04 01:01:58 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 28 17:32:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	199377	20.00	ng	-0.03
21) Naphthalene-d8	7.96	136	730234	20.00	ng	-0.04
38) Acenaphthene-d10	9.70	164	350075	20.00	ng	-0.05
63) Phenanthrene-d10	11.19	188	563946	20.00	ng	-0.04
75) Chrysene-d12	13.82	240	380210	20.00	ng	-0.05
86) Perylene-d12	15.20	264	344851	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1460069	122.28	ng	-0.03
7) Phenol-d6	6.36	99	2112937	144.94	ng	-0.03
23) Nitrobenzene-d5	7.25	82	1460800	111.24	ng	-0.04
41) 2,4,6-Tribromophenol	10.50	330	468006	161.54	ng	-0.04
44) 2-Fluorobiphenyl	9.04	172	1968652	120.63	ng	-0.04
78) Terphenyl-d14	12.78	244	974245	62.14	ng	-0.04
Target Compounds						
2) 1,4-Dioxane	2.28	88	117150	19.93	ng	# 74
4) n-Nitrosodimethylamine	2.81	42	58891	7.61	ng	# 82
10) Phenol	6.37	94	117185	7.05	ng	# 62
15) Benzyl Alcohol	6.84	79	158236	13.97	ng	96
20) 3+4-Methylphenols	7.11	107	60206	4.62	ng	# 68
31) Naphthalene	7.98	128	110832	3.32	ng	99
32) Benzoic acid	8.12	122	11980725m	1411.94	ng	
48) Acenaphthylene	9.46	152	878749	30.10	ng	# 78
59) Diethylphthalate	10.14	149	171045	7.87	ng	99

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

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