

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111546.D
 Acq On : 17 Dec 2018 18:36
 Operator : JU/SJ
 Sample : J6438-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:49:19 PM

Quant Time: Dec 18 06:11:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	143950	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	575287	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	407692	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	688465	20.00	ng	0.00
76) Chrysene-d12	13.95	240	454070	20.00	ng	0.00
87) Perylene-d12	15.39	264	440400	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	653514	75.55	ng	0.00
7) Phenol-d6	6.45	99	462244	40.17	ng	0.00
23) Nitrobenzene-d5	7.36	82	922010	95.44	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	486410	133.19	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1909542	72.35	ng	0.00
79) Terphenyl-d14	12.90	244	2005753	103.73	ng	0.00
Target Compounds						
9) Benzaldehyde	6.33	77	31648	4.535	ng	98
10) Phenol	6.46	94	70128	5.471	ng	92
15) Benzyl Alcohol	6.94	79	96710	11.046	ng	90
17) 2-Methylphenol	7.07	107	28206	3.391	ng	93
20) 3+4-Methylphenols	7.22	107	286177	27.426	ng	# 63
27) 2,4-Dimethylphenol	7.74	122	134504	18.153	ng	98
31) Naphthalene	8.09	128	172666	6.417	ng	95
32) Benzoic acid	7.96	122	419169m	65.467	ng	
50) Dimethylphthalate	9.55	163	77479	2.639	ng	99

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

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