

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111545.D
 Acq On : 17 Dec 2018 18:08
 Operator : JU/SJ
 Sample : J6438-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:49:17 PM

Quant Time: Dec 18 06:07:04 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	125213	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	500461	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	341013	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	553498	20.00	ng	0.00
76) Chrysene-d12	13.94	240	358313	20.00	ng	0.00
87) Perylene-d12	15.39	264	376279	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	598756	79.57	ng	0.01
7) Phenol-d6	6.45	99	398402	39.80	ng	0.00
23) Nitrobenzene-d5	7.35	82	772956	91.97	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	369790	121.05	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1448158	65.60	ng	0.00
79) Terphenyl-d14	12.90	244	1457033	95.49	ng	0.00
Target Compounds						
9) Benzaldehyde	6.33	77	23384	3.852	ng	97
10) Phenol	6.46	94	50724	4.549	ng	96
15) Benzyl Alcohol	6.94	79	285932	37.545	ng	93
20) 3+4-Methylphenols	7.22	107	64913	7.152	ng	# 58
27) 2,4-Dimethylphenol	7.75	122	24446	3.793	ng	97
32) Benzoic acid	7.96	122	435076m	78.112	ng	
50) Dimethylphthalate	9.54	163	84146	3.426	ng	# 93
66) n-Nitrosodiphenylamine	10.48	169	62954	3.298	ng	94

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

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