

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091585.D
 Acq On : 14 Dec 2016 14:51
 Operator : UM/SJ
 Sample : H5961-01DL 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW26S-GWDL

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:39:42 PM

Quant Time: Dec 15 14:07:13 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	265408	20.00	ng	-0.02
21) Naphthalene-d8	8.09	136	1051144	20.00	ng	-0.01
38) Acenaphthene-d10	9.84	164	481695	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	681005	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	531708	20.00	ng	-0.01
86) Perylene-d12	15.37	264	570274	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	173092	10.99	ng	-0.02
7) Phenol-d6	6.43	99	147663	7.52	ng	-0.02
23) Nitrobenzene-d5	7.36	82	286254	15.20	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	94685	23.67	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	536451	19.24	ng	-0.02
78) Terphenyl-d14	12.91	244	275196	11.74	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	2.36	88	25634	3.08	ng	# 80
6) Aniline	6.45	93	55868	2.17	ng	# 92
10) Phenol	6.45	94	83678	3.64	ng	85
17) 2-Methylphenol	7.06	107	57456m	3.91	ng	
20) 3+4-Methylphenols	7.22	107	224826	11.85	ng	# 88
27) 2,4-Dimethylphenol	7.76	122	222715	14.44	ng	90
31) Naphthalene	8.12	128	7283540	148.93	ng	94
37) 2-Methylnaphthalene	8.80	142	277025	8.75	ng	97
45) 1,1'-Biphenyl	9.26	154	436312	12.51	ng	96
51) Acenaphthene	9.87	154	820049	28.67	ng	97
54) Dibenzofuran	10.04	168	792455	22.39	ng	100
57) Fluorene	10.38	166	531404	18.03	ng	98
70) Phenanthrene	11.34	178	845473	25.07	ng	100
71) Anthracene	11.40	178	98092	2.70	ng	99
72) Carbazole	11.55	167	390345	12.24	ng	98
74) Fluoranthene	12.53	202	79099	2.24	ng	91

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

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