

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081517\
 Data File : BF097689.D
 Acq On : 15 Aug 2017 14:21
 Operator : SJ/JU
 Sample : I4705-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-B-COMP

Manual Integrations
 APPROVED

Sohil
 8/16/2017 6:45:14 PM

Quant Time: Aug 15 15:28:29 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	130264	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	525004	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	237630	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	467775	20.00	ng	0.00
75) Chrysene-d12	13.93	240	337766	20.00	ng	0.00
86) Perylene-d12	15.36	264	260038	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	856803	100.05	ng	0.00
7) Phenol-d6	6.41	99	1173993	100.89	ng	0.00
23) Nitrobenzene-d5	7.35	82	703111	72.75	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	221974	107.66	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	975940	60.34	ng	0.00
78) Terphenyl-d14	12.89	244	905582	55.91	ng	0.00
Target Compounds						
10) Phenol	6.42	94	34870	2.562	ng	Qvalue # 75
49) Dimethylphthalate	9.53	163	249437	14.360	ng	100
74) Fluoranthene	12.50	202	61993	2.383	ng	98
82) Chrysene	13.96	228	41998	2.086	ng	96
85) Indeno(1,2,3-cd)pyrene	16.74	276	38328m	2.587	ng	
87) Benzo(b)fluoranthene	14.95	252	131599m	8.029	ng	
89) Benzo(a)pyrene	15.30	252	65199	4.493	ng	99
91) Benzo(g,h,i)perylene	17.16	276	31650	2.568	ng	# 90

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

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