

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072017\
 Data File : BF096930.D
 Acq On : 21 Jul 2017 4:04
 Operator : SJ/JU
 Sample : I4253-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-O-P-6-7-WSW

Manual Integrations
 APPROVED

Sohil
 7/21/2017 1:06:09 PM

Quant Time: Jul 21 08:45:41 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 13:33:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.57	152	113977	20.00	ng	-0.03
21) Naphthalene-d8	7.85	136	455534	20.00	ng	-0.04
38) Acenaphthene-d10	9.60	164	173839	20.00	ng	-0.04
63) Phenanthrene-d10	11.07	188	251104	20.00	ng	-0.03
75) Chrysene-d12	13.70	240	187067	20.00	ng	-0.03
86) Perylene-d12	15.07	264	152468	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	958253	126.41	ng	-0.02
7) Phenol-d6	6.23	99	1201032	132.03	ng	-0.02
23) Nitrobenzene-d5	7.14	82	723169	98.33	ng	-0.03
41) 2,4,6-Tribromophenol	10.39	330	181794	122.51	ng	-0.04
44) 2-Fluorobiphenyl	8.93	172	954524	86.75	ng	-0.03
78) Terphenyl-d14	12.66	244	575758	53.38	ng	-0.03
Target Compounds						Qvalue
10) Phenol	6.24	94	33681	3.275	ng	96
31) Naphthalene	7.87	128	49705	2.303	ng	99
49) Dimethylphthalate	9.33	163	251059	19.180	ng	99
51) Acenaphthene	9.63	154	22111	2.140	ng	94
57) Fluorene	10.14	166	22899	2.140	ng	# 97
70) Phenanthrene	11.10	178	303191	22.206	ng	98
71) Anthracene	11.15	178	96380	6.982	ng	98
72) Carbazole	11.30	167	37771	2.825	ng	96
74) Fluoranthene	12.29	202	537396	41.121	ng	98
77) Pyrene	12.51	202	523654	30.844	ng	98
80) Benzo(a)anthracene	13.70	228	315036	27.029	ng	96
82) Chrysene	13.73	228	285037	24.841	ng	97
85) Indeno(1,2,3-cd)pyrene	16.35	276	166877	18.132	ng	93
87) Benzo(b)fluoranthene	14.70	252	441526m	48.015	ng	
88) Benzo(k)fluoranthene	14.73	252	105555m	12.122	ng	
89) Benzo(a)pyrene	15.02	252	276777	32.817	ng	97
90) Dibenzo(a,h)anthracene	16.50	278	37596	4.845	ng	93
91) Benzo(g,h,i)perylene	16.73	276	150553	18.116	ng	98

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

