

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092217\
 Data File : BF098689.D
 Acq On : 22 Sep 2017 10:16
 Operator : SJ/JU
 Sample : I5311-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-091917

Manual Integrations
 APPROVED

Sohil
 9/25/2017 6:20:56 PM

Quant Time: Sep 22 12:58:24 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 11:54:28 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	144043	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	562662	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	215796	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	305537	20.00	ng	0.00
75) Chrysene-d12	14.09	240	282816	20.00	ng	0.00
86) Perylene-d12	15.58	264	325554	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	725715	76.51	ng	0.00
7) Phenol-d6	6.55	99	1031305	89.78	ng	0.00
23) Nitrobenzene-d5	7.49	82	645660	77.84	ng	0.00
41) 2,4,6-Tribromophenol	10.75	330	87690	44.43	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	1059648	78.14	ng	0.00
78) Terphenyl-d14	13.03	244	739033	56.11	ng	0.00
Target Compounds						
10) Phenol	6.55	94	30539	2.409	ng	86
49) Dimethylphthalate	9.67	163	86039	5.317	ng	99
70) Phenanthrene	11.47	178	64236	3.876	ng	96
74) Fluoranthene	12.66	202	174252	10.549	ng	98
77) Pyrene	12.89	202	184849	8.804	ng	98
80) Benzo(a)anthracene	14.07	228	123654	7.259	ng	# 78
82) Chrysene	14.11	228	127118	7.727	ng	97
85) Indeno(1,2,3-cd)pyrene	17.09	276	101331	7.940	ng	96
87) Benzo(b)fluoranthene	15.13	252	225835m	11.961	ng	
88) Benzo(k)fluoranthene	15.16	252	66969m	3.549	ng	
89) Benzo(a)pyrene	15.52	252	151837	8.855	ng	# 93
91) Benzo(a,h,i)perylene	17.54	276	97590	6.695	ng	98

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

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