

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112117\
 Data File : BF100793.D
 Acq On : 21 Nov 2017 21:00
 Operator : SJ/JU
 Sample : I6499-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PG1-1-E(7-8)

Manual Integrations
 APPROVED

Sohil
 11/22/2017 5:04:28 PM

Quant Time: Nov 22 06:47:18 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 15:22:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	73543	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	281601	20.00	ng	-0.02
38) Acenaphthene-d10	9.91	164	118952	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	185284	20.00	ng	-0.01
75) Chrysene-d12	14.03	240	156875	20.00	ng	-0.01
86) Perylene-d12	15.50	264	167519	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	479848	104.59	ng	0.00
7) Phenol-d6	6.50	99	616649	109.00	ng	0.00
23) Nitrobenzene-d5	7.43	82	372403	87.43	ng	-0.02
41) 2,4,6-Tribromophenol	10.70	330	136518	112.63	ng	-0.01
44) 2-Fluorobiphenyl	9.23	172	740888	94.84	ng	-0.01
78) Terphenyl-d14	12.98	244	499182	73.11	ng	0.00
Target Compounds						
10) Phenol	6.51	94	28580	4.812	ng	90
49) Dimethylphthalate	9.62	163	68372	7.528	ng	99
70) Phenanthrene	11.42	178	39541	4.053	ng	98
74) Fluoranthene	12.60	202	73419	7.101	ng	97
77) Pyrene	12.83	202	67529	6.039	ng	97
80) Benzo(a)anthracene	14.02	228	49503	5.240	ng	96
82) Chrysene	14.06	228	41671	4.555	ng	97
85) Indeno(1,2,3-cd)pyrene	16.98	276	23372	3.159	ng	93
87) Benzo(b)fluoranthene	15.07	252	63744m	6.423	ng	
89) Benzo(a)pyrene	15.44	252	42094	4.784	ng	97
91) Benzo(g,h,i)perylene	17.43	276	21229	3.014	ng	98

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

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