

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103464.D
 Acq On : 6 Mar 2018 19:16
 Operator : SJ/JU
 Sample : J1722-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-18

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:07 PM

Quant Time: Mar 07 00:56:40 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 11:52:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	150767	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	662412	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	282638	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	438240	20.00	ng	0.00
75) Chrysene-d12	14.06	240	279374	20.00	ng	0.00
86) Perylene-d12	15.58	264	234968	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	868025	87.08	ng	0.00
7) Phenol-d6	6.51	99	1071616	87.85	ng	0.00
23) Nitrobenzene-d5	7.44	82	642429	55.39	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	130150	54.40	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	871053	47.49	ng	0.00
78) Terphenyl-d14	12.99	244	565371	48.65	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	53834	2.374	ng	Qvalue 100
70) Phenanthrene	11.43	178	246883	10.237	ng	98
71) Anthracene	11.49	178	80456	3.253	ng	99
74) Fluoranthene	12.63	202	313975	12.955	ng	99
77) Pyrene	12.86	202	283757	13.027	ng	99
80) Benzo(a)anthracene	14.05	228	122247	6.744	ng	98
82) Chrysene	14.09	228	110564	6.282	ng	98
85) Indeno(1,2,3-cd)pyrene	17.12	276	40944	2.938	ng	97
87) Benzo(b)fluoranthene	15.13	252	115396m	7.470	ng	
88) Benzo(k)fluoranthene	15.16	252	41200m	2.832	ng	
89) Benzo(a)pyrene	15.52	252	83007	6.017	ng	# 92
91) Benzo(a,h,i)perylene	17.60	276	38235	3.670	ng	92

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

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