

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103469.D
 Acq On : 6 Mar 2018 21:29
 Operator : SJ/JU
 Sample : J1722-13 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-9

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:24 PM

Quant Time: Mar 07 01:15:55 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	125705	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	536349	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	215409	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	331433	20.00	ng	0.00
75) Chrysene-d12	14.06	240	217302	20.00	ng	0.00
86) Perylene-d12	15.57	264	165847	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	464037	55.83	ng	0.00
7) Phenol-d6	6.51	99	626456	61.60	ng	0.00
23) Nitrobenzene-d5	7.44	82	360507	38.39	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	63045	34.58	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	508781	34.70	ng	0.00
78) Terphenyl-d14	12.99	244	344669	38.13	ng	-0.01
Target Compounds						
70) Phenanthrene	11.43	178	100329	5.501	ng	98
74) Fluoranthene	12.63	202	169467	9.246	ng	99
77) Pyrene	12.86	202	168209	9.928	ng	99
80) Benzo(a)anthracene	14.05	228	72146	5.117	ng	98
82) Chrysene	14.09	228	65669	4.797	ng	97
83) Bis(2-ethylhexyl)phthalate	14.01	149	47318	3.917	ng	98
85) Indeno(1,2,3-cd)pyrene	17.12	276	23619	2.179	ng	99
87) Benzo(b)fluoranthene	15.13	252	63994m	5.869	ng	
88) Benzo(k)fluoranthene	15.16	252	22192m	2.161	ng	
89) Benzo(a)pyrene	15.52	252	42400	4.354	ng	# 88
91) Benzo(g,h,i)perylene	17.59	276	22855	3.108	ng	# 84

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

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