

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
 Data File : BF103788.D
 Acq On : 20 Mar 2018 6:50
 Operator : JU/SJ
 Sample : J1973-02 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2

Manual Integrations
 APPROVED

Sohil
 3/20/2018 6:27:50 PM

Quant Time: Mar 20 08:06:29 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	141691	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	562459	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	253213	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	411128	20.00	ng	0.00
75) Chrysene-d12	14.16	240	292875	20.00	ng	0.00
86) Perylene-d12	15.69	264	260055	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	187808	20.16	ng	0.00
7) Phenol-d6	6.58	99	256187	22.48	ng	0.00
23) Nitrobenzene-d5	7.53	82	143600	17.80	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	27678	12.71	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	192570	12.13	ng	0.00
78) Terphenyl-d14	13.09	244	149096	11.82	ng	0.00
Target Compounds						
70) Phenanthrene	11.53	178	103834	4.617	ng	98
71) Anthracene	11.58	178	54178	2.372	ng	99
74) Fluoranthene	12.72	202	136687	5.899	ng	98
77) Pyrene	12.95	202	149942	6.971	ng	98
80) Benzo(a)anthracene	14.14	228	79552m	4.291	ng	
82) Chrysene	14.18	228	85074m	4.814	ng	
85) Indeno(1,2,3-cd)pyrene	17.26	276	31049	2.012	ng	93
87) Benzo(b)fluoranthene	15.23	252	73122m	4.451	ng	
89) Benzo(a)pyrene	15.62	252	66209	4.443	ng	# 94
91) Benzo(g,h,i)perylene	17.73	276	30497	2.429	ng	96

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

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