

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112817\
 Data File : BF100935.D
 Acq On : 28 Nov 2017 15:48
 Operator : SJ/JU
 Sample : I6579-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HAR-4-E(0-1)

Manual Integrations
 APPROVED

Sohil
 11/29/2017 4:28:32 PM

Quant Time: Nov 29 06:41:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 15:21:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	66449	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	256499	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	108860	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	172684	20.00	ng	0.00
75) Chrysene-d12	14.02	240	152225	20.00	ng	0.00
86) Perylene-d12	15.49	264	158662	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	394599	95.19	ng	0.00
7) Phenol-d6	6.49	99	479882	93.88	ng	0.00
23) Nitrobenzene-d5	7.42	82	298276	76.88	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	111774	100.76	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	554406	77.55	ng	0.00
78) Terphenyl-d14	12.96	244	406034	61.29	ng	0.00
Target Compounds						
10) Phenol	6.50	94	11770	2.193	ng	# 57
49) Dimethylphthalate	9.60	163	35168	4.231	ng	# 99
70) Phenanthrene	11.40	178	22703	2.497	ng	95
74) Fluoranthene	12.59	202	71211	7.390	ng	95
77) Pyrene	12.82	202	73147	6.741	ng	98
80) Benzo(a)anthracene	14.01	228	48317	5.271	ng	97
82) Chrysene	14.04	228	40836	4.600	ng	96
85) Indeno(1,2,3-cd)pyrene	16.96	276	26449	3.684	ng	# 91
87) Benzo(b)fluoranthene	15.06	252	63086m	6.711	ng	
88) Benzo(k)fluoranthene	15.09	252	18543m	2.088	ng	
89) Benzo(a)pyrene	15.43	252	46783	5.614	ng	99
91) Benzo(a,h,i)perylene	17.41	276	25036	3.753	ng	# 90

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

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