

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
 Data File : BF098391.D
 Acq On : 9 Sep 2017 17:43
 Operator : SJ/JU
 Sample : I5133-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-204

Manual Integrations
 APPROVED

Sohil
 9/11/2017 2:42:45 PM

Quant Time: Sep 11 11:59:46 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	114177	20.00	ng	-0.03
21) Naphthalene-d8	7.96	136	439693	20.00	ng	-0.03
38) Acenaphthene-d10	9.70	164	186585	20.00	ng	-0.04
63) Phenanthrene-d10	11.19	188	297605	20.00	ng	-0.04
75) Chrysene-d12	13.82	240	223114	20.00	ng	-0.04
86) Perylene-d12	15.23	264	218046	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	406778	54.19	ng	-0.02
7) Phenol-d6	6.33	99	567450	55.64	ng	-0.02
23) Nitrobenzene-d5	7.24	82	330436	40.83	ng	-0.04
41) 2,4,6-Tribromophenol	10.49	330	87276	53.91	ng	-0.04
44) 2-Fluorobiphenyl	9.03	172	524834	41.33	ng	-0.04
78) Terphenyl-d14	12.77	244	330552	30.90	ng	-0.04
Target Compounds						
49) Dimethylphthalate	9.43	163	69621	5.105	ng	Qvalue # 89
70) Phenanthrene	11.21	178	344889	21.748	ng	99
71) Anthracene	11.26	178	100957	6.277	ng	97
74) Fluoranthene	12.40	202	854400	51.617	ng	99
77) Pyrene	12.63	202	772826	42.351	ng	99
80) Benzo(a)anthracene	13.82	228	418309	31.282	ng	98
82) Chrysene	13.85	228	378668	28.467	ng	98
85) Indeno(1,2,3-cd)pyrene	16.57	276	176751	18.062	ng	# 92
87) Benzo(b)fluoranthene	14.83	252	508301m	36.983	ng	
88) Benzo(k)fluoranthene	14.86	252	145169m	11.242	ng	
89) Benzo(a)pyrene	15.17	252	355880	29.249	ng	98
90) Dibenzo(a,h)anthracene	16.57	278	46542m	4.699	ng	
91) Benzo(a,h,i)perylene	16.98	276	171595	16.602	ng	93

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

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