

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112817\
 Data File : BF100934.D
 Acq On : 28 Nov 2017 15:22
 Operator : SJ/JU
 Sample : I6578-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BBR-2-E(7-8)

Manual Integrations
 APPROVED

Sohil
 11/29/2017 4:28:30 PM

Quant Time: Nov 29 06:32:56 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	68380	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	245402	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	101353	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	155854	20.00	ng	0.00
75) Chrysene-d12	14.02	240	150587	20.00	ng	0.00
86) Perylene-d12	15.49	264	162372	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	382987	89.78	ng	0.00
7) Phenol-d6	6.49	99	446521	84.89	ng	0.00
23) Nitrobenzene-d5	7.42	82	271757	73.21	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	97884	94.78	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	569027	85.49	ng	0.00
78) Terphenyl-d14	12.96	244	406379	62.01	ng	0.00
Target Compounds						
10) Phenol	6.50	94	11331	2.052	ng	77
49) Dimethylphthalate	9.60	163	23600	3.050	ng	# 99
70) Phenanthrene	11.40	178	39249	4.783	ng	97
74) Fluoranthene	12.59	202	87347	10.044	ng	95
77) Pyrene	12.82	202	88945	8.286	ng	98
80) Benzo(a)anthracene	14.01	228	48274m	5.323	ng	
82) Chrysene	14.04	228	55469m	6.317	ng	
85) Indeno(1,2,3-cd)pyrene	16.96	276	35055	4.935	ng	# 88
87) Benzo(b)fluoranthene	15.06	252	78580m	8.169	ng	
88) Benzo(k)fluoranthene	15.09	252	22132m	2.435	ng	
89) Benzo(a)pyrene	15.43	252	53695	6.296	ng	97
91) Benzo(a,h,i)perylene	17.41	276	31989	4.686	ng	# 93

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

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