

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112817\
 Data File : BF100932.D
 Acq On : 28 Nov 2017 14:28
 Operator : SJ/JU
 Sample : I6578-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BBR-2-E(0-1)

Manual Integrations
 APPROVED

Sohil
 11/29/2017 4:28:27 PM

Quant Time: Nov 29 06:20:14 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	67617	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	255874	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	111971	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	177044	20.00	ng	0.00
75) Chrysene-d12	14.02	240	146006	20.00	ng	0.00
86) Perylene-d12	15.50	264	151195	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	475218	112.66	ng	0.00
7) Phenol-d6	6.49	99	602090	115.75	ng	0.00
23) Nitrobenzene-d5	7.43	82	369158	95.38	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	144220	126.40	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	780538	106.15	ng	0.00
78) Terphenyl-d14	12.97	244	552758	86.99	ng	0.00
Target Compounds						
10) Phenol	6.50	94	11502	2.106	ng	Qvalue 97
49) Dimethylphthalate	9.60	163	37711	4.411	ng	100
74) Fluoranthene	12.59	202	24818	2.512	ng	96
77) Pyrene	12.82	202	25450	2.445	ng	99
82) Chrysene	14.04	228	20297m	2.384	ng	
87) Benzo(b)fluoranthene	15.06	252	23067m	2.575	ng	
89) Benzo(a)pyrene	15.43	252	17488	2.202	ng	# 82
91) Benzo(g,h,i)perylene	17.43	276	15190	2.389	ng	# 92

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

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