

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041117\
 Data File : BF094381.D
 Acq On : 12 Apr 2017 6:58
 Operator : SJ/MA
 Sample : I2598-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-64-S1(0-2)

Manual Integrations
 APPROVED

mohammad
 4/12/2017 5:12:23 PM

Quant Time: Apr 12 07:58:00 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 06 13:20:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.83	152	137753	20.00	ng	-0.03
21) Naphthalene-d8	7.33	136	517506	20.00	ng	-0.04
38) Acenaphthene-d10	9.47	164	204618	20.00	ng	-0.03
63) Phenanthrene-d10	11.29	188	331238	20.00	ng	-0.03
75) Chrysene-d12	14.54	240	293885	20.00	ng	-0.03
86) Perylene-d12	16.17	264	212992	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	4.38	112	981046	120.29	ng	0.00
7) Phenol-d6	5.46	99	1338710	132.68	ng	0.00
23) Nitrobenzene-d5	6.48	82	844845	93.17	ng	-0.03
41) 2,4,6-Tribromophenol	10.45	330	142129	87.90	ng	-0.03
44) 2-Fluorobiphenyl	8.66	172	1324783	98.75	ng	-0.03
78) Terphenyl-d14	13.27	244	905520	69.06	ng	-0.02
Target Compounds						
10) Phenol	5.47	94	81584	7.11	ng	Qvalue 94
49) Dimethylphthalate	9.15	163	172416	11.85	ng	98
74) Fluoranthene	12.78	202	64238	3.40	ng	99
77) Pyrene	13.06	202	67075	3.14	ng	98
80) Benzo(a)anthracene	14.53	228	35113	2.05	ng	98
82) Chrysene	14.57	228	43438	2.73	ng	98
87) Benzo(b)fluoranthene	15.77	252	55293m	4.24	ng	
89) Benzo(a)pyrene	16.11	252	34324	2.85	ng	# 95
91) Benzo(g,h,i)perylene	17.87	276	20857	2.03	ng	97

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

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