

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080516\
 Data File : BF089629.D
 Acq On : 5 Aug 2016 14:21
 Operator : UM/SJ
 Sample : H4292-02
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS6-0-2IN

Manual Integrations
 APPROVED

umangi
 8/8/2016 9:09:34 AM

Quant Time: Aug 05 23:24:47 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 01:39:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	50369	20.00	ng	-0.01
21) Naphthalene-d8	9.46	136	215320	20.00	ng	-0.01
38) Acenaphthene-d10	12.28	164	97899	20.00	ng	-0.01
63) Phenanthrene-d10	14.69	188	172036	20.00	ng	-0.01
75) Chrysene-d12	18.39	240	100886	20.00	ng	0.00
86) Perylene-d12	20.05	264	101160	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	292728	97.41	ng	0.02
7) Phenol-d6	6.98	99	389312	95.50	ng	0.00
23) Nitrobenzene-d5	8.35	82	248250	63.77	ng	-0.01
41) 2,4,6-Tribromophenol	13.58	330	73513	90.99	ng	-0.01
44) 2-Fluorobiphenyl	11.24	172	455044	60.41	ng	-0.01
78) Terphenyl-d14	17.05	244	309218	60.51	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	11.94	163	81824	10.66	ng	99
70) Phenanthrene	14.72	178	67668	7.23	ng	100
71) Anthracene	14.82	178	24426	2.44	ng	96
74) Fluoranthene	16.49	202	143838	14.96	ng	97
77) Pyrene	16.79	202	121926	13.67	ng	96
80) Benzo(a)anthracene	18.38	228	62612m	10.80	ng	
82) Chrysene	18.42	228	53034	8.72	ng	97
85) Indeno(1,2,3-cd)pyrene	21.23	276	43146	9.34	ng	94
87) Benzo(b)fluoranthene	19.65	252	71341m	11.47	ng	
88) Benzo(k)fluoranthene	19.66	252	38133m	6.08	ng	
89) Benzo(a)pyrene	19.99	252	60392	10.37	ng	# 91
91) Benzo(g,h,i)perylene	21.57	276	40312	7.24	ng	98

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

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