

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110416\
 Data File : BF091034.D
 Acq On : 4 Nov 2016 21:35
 Operator : UM/SJ
 Sample : H5526-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-98-4.5-5.0

Manual Integrations
 APPROVED

sohil
 11/7/2016 4:39:46 PM

Quant Time: Nov 04 23:58:34 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 11:03:12 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	176031	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	682441	20.00	ng	-0.01
38) Acenaphthene-d10	9.74	164	336226	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	452962	20.00	ng	-0.01
75) Chrysene-d12	13.85	240	329348	20.00	ng	-0.01
86) Perylene-d12	15.24	264	294518	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	1684331	158.03	ng	0.02
7) Phenol-d6	6.36	99	2170202	150.01	ng	0.01
23) Nitrobenzene-d5	7.28	82	1454673	116.28	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	410771	135.14	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1828990	93.65	ng	0.00
78) Terphenyl-d14	12.81	244	1328273	100.64	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.47	163	587253	26.64	ng	99
74) Fluoranthene	12.43	202	172249	6.90	ng	91
77) Pyrene	12.66	202	141450	6.56	ng	95
80) Benzo(a)anthracene	13.85	228	96662	5.17	ng	99
82) Chrysene	13.88	228	87427m	4.74	ng	
85) Indeno(1,2,3-cd)pyrene	16.55	276	58560	3.83	ng	93
87) Benzo(b)fluoranthene	14.85	252	133425m	6.68	ng	
88) Benzo(k)fluoranthene	14.87	252	52323m	2.99	ng	
89) Benzo(a)pyrene	15.19	252	78148	4.49	ng	# 97
91) Benzo(g,h,i)perylene	16.93	276	55030	4.05	ng	99

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

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