

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110416\
 Data File : BF091030.D
 Acq On : 4 Nov 2016 19:45
 Operator : UM/SJ
 Sample : H5526-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-101-1.0-1.5

Manual Integrations
 APPROVED

sohil
 11/7/2016 4:39:58 PM

Quant Time: Nov 07 04:06:17 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	224453	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	932812	20.00	ng	-0.01
38) Acenaphthene-d10	9.74	164	488328	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	729111	20.00	ng	-0.01
75) Chrysene-d12	13.86	240	514927	20.00	ng	0.00
86) Perylene-d12	15.24	264	416194	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1413523	104.01	ng	0.01
7) Phenol-d6	6.35	99	1673979	90.75	ng	0.00
23) Nitrobenzene-d5	7.28	82	1252566	73.25	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	438680	99.37	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1799066	63.43	ng	0.00
78) Terphenyl-d14	12.81	244	1334095	64.65	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.47	163	576171	17.99	ng	98
70) Phenanthrene	11.24	178	1221433	31.51	ng	99
71) Anthracene	11.30	178	235303	5.71	ng	99
72) Carbazole	11.46	167	152802	4.11	ng	97
74) Fluoranthene	12.43	202	1884537	46.88	ng	98
77) Pyrene	12.66	202	1536678	45.57	ng	99
80) Benzo(a)anthracene	13.85	228	716637m	24.51	ng	
82) Chrysene	13.88	228	705347m	24.47	ng	
85) Indeno(1,2,3-cd)pyrene	16.56	276	397435	16.62	ng	93
87) Benzo(b)fluoranthene	14.87	252	927681m	32.86	ng	
88) Benzo(k)fluoranthene	14.88	252	192448m	7.77	ng	
89) Benzo(a)pyrene	15.19	252	535747	21.80	ng	# 94
90) Dibenzo(a,h)anthracene	16.56	278	90353m	4.25	ng	
91) Benzo(g,h,i)perylene	16.95	276	364763	18.99	ng	99

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

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