

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
 Data File : BF086436.D
 Acq On : 27 Apr 2016 21:03
 Operator : UM/SJ
 Sample : H2724-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TH-B4(0-5)-C

Manual Integrations
 APPROVED

sohil
 4/28/2016 1:40:08 PM

Quant Time: Apr 28 01:45:11 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	126995	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	522644	20.00	ng	-0.01
38) Acenaphthene-d10	9.84	164	211240	20.00	ng	-0.01
63) Phenanthrene-d10	11.32	188	324339	20.00	ng	0.00
75) Chrysene-d12	13.95	240	271436	20.00	ng	-0.01
86) Perylene-d12	15.36	264	268867	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	831946	113.00	ng	0.01
7) Phenol-d6	6.44	99	1074930	104.61	ng	0.00
23) Nitrobenzene-d5	7.37	82	638321	65.83	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	177514	79.69	ng	-0.01
44) 2-Fluorobiphenyl	9.17	172	1084192	91.23	ng	0.00
78) Terphenyl-d14	12.91	244	659927	53.56	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.56	163	146648	9.22	ng	99
70) Phenanthrene	11.34	178	53328	3.13	ng	100
74) Fluoranthene	12.53	202	121758	7.09	ng	93
77) Pyrene	12.75	202	105058	5.37	ng	97
80) Benzo(a)anthracene	13.94	228	65555	4.53	ng	96
82) Chrysene	13.97	228	62443m	3.97	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	36806	3.16	ng	# 92
87) Benzo(b)fluoranthene	14.96	252	106835m	6.75	ng	
88) Benzo(k)fluoranthene	14.98	252	35446m	2.15	ng	
89) Benzo(a)pyrene	15.30	252	63898	4.28	ng	# 93
91) Benzo(g,h,i)perylene	17.12	276	36175	2.95	ng	97

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

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